

4/12/05

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NEWS 3 FEB 25 CA/CAPLUS - Russian Agency for Patents and Trademarks (ROSPATENT) added to list of core patent offices covered  
NEWS 4 FEB 28 PATDPAFULL - New display fields provide for legal status data from INPADOC  
NEWS 5 FEB 28 BABS - Current-awareness alerts (SDIs) available  
NEWS 6 FEB 28 MEDLINE/LMEDLINE reloaded  
NEWS 7 MAR 02 GBFULL: New full-text patent database on STN  
NEWS 8 MAR 03 REGISTRY/ZREGISTRY - Sequence annotations enhanced  
NEWS 9 MAR 03 MEDLINE file segment of TOXCENTER reloaded  
NEWS 10 MAR 22 KOREPAT now updated monthly; patent information enhanced  
NEWS 11 MAR 22 Original IDE display format returns to REGISTRY/ZREGISTRY  
NEWS 12 MAR 22 PATDPASPC - New patent database available  
NEWS 13 MAR 22 REGISTRY/ZREGISTRY enhanced with experimental property tags  
NEWS 14 APR 04 EPFULL enhanced with additional patent information and new fields  
NEWS 15 APR 04 EMBASE - Database reloaded and enhanced

NEWS EXPRESS JANUARY 10 CURRENT WINDOWS VERSION IS V7.01a, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 10 JANUARY 2005

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FILE 'HOME' ENTERED AT 14:31:03 ON 13 APR 2005

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=> file registry			
COST IN U.S. DOLLARS		SINCE FILE	TOTAL
		ENTRY	SESSION
FULL ESTIMATED COST		0.21	0.21

FILE 'REGISTRY' ENTERED AT 14:31:11 ON 13 APR 2005  
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STRUCTURE FILE UPDATES: 12 APR 2005 HIGHEST RN 848391-87-7  
DICTIONARY FILE UPDATES: 12 APR 2005 HIGHEST RN 848391-87-7

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when  
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\*\*\*\*\*  
\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more  
information enter HELP PROP at an arrow prompt in the file or refer  
to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

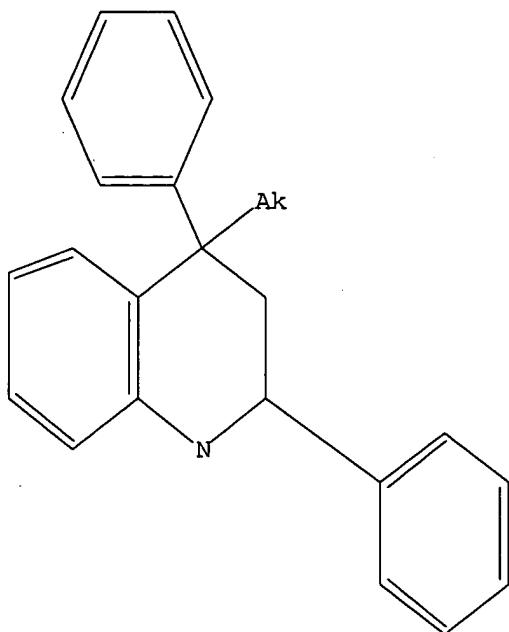
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L1 STRUCTURE UPLOADED

=> d 11  
L1 HAS NO ANSWERS  
L1 STR

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G1 O,S,CH2

G2 C,N

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SEARCH INITIATED 14:31:34 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 4738 TO ITERATE

21.1% PROCESSED 1000 ITERATIONS 2 ANSWERS  
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 90633 TO 98887  
PROJECTED ANSWERS: 5 TO 373

L2 2 SEA SSS SAM L1

=> s 11 ful  
FULL SEARCH INITIATED 14:31:40 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 95702 TO ITERATE

100.0% PROCESSED 95702 ITERATIONS 119 ANSWERS  
SEARCH TIME: 00.00.03

L3 119 SEA SSS FUL L1

=> file caplus  
COST IN U.S. DOLLARS SINCE FILE TOTAL  
ENTRY SESSION

10796396

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FULL ESTIMATED COST

161.33

161.54

FILE 'CAPLUS' ENTERED AT 14:31:47 ON 13 APR 2005  
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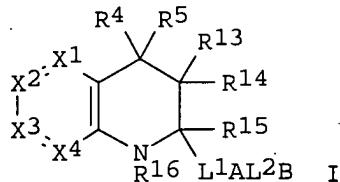
FILE COVERS 1907 - 13 Apr 2005 VOL 142 ISS 16  
FILE LAST UPDATED: 12 Apr 2005 (20050412/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13  
L4 10 L3

=> d abs bib hitstr 1-10

L4 ANSWER 1 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
GI



AB Title compds. [I; L1 = bond, CH<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>, CH<sub>2</sub>O, CH<sub>2</sub>CO, etc.; L2 = bond, O, CO, CO<sub>2</sub>, S, SO, SO<sub>2</sub>, CONR<sub>8</sub>, SO<sub>2</sub>NR<sub>8</sub>, etc.; A = (substituted) carbocyclene, heterocyclene; B = (substituted) alkyl, alkenyl, alkynyl, carbocyclyl, heterocyclyl; X<sub>1</sub>-X<sub>4</sub> = CR<sub>1</sub>, CR<sub>2</sub>, N, etc.; R<sub>1</sub> = H, F, Cl, Br, iodo, OCF<sub>3</sub>, CF<sub>3</sub>, cyano, NH<sub>2</sub>, alkylamino, dialkylamino, CONH<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>, etc.; R<sub>2</sub> = H, F, Cl, Br, iodo, OCF<sub>3</sub>, CF<sub>3</sub>, cyano, NO<sub>2</sub>, amino, aminocarbonyl, (substituted) alkyl, alkenyl, alkynyl, carbocyclyl, heterocyclyl, etc.; R<sub>4</sub> = H, F, haloalkyl, (substituted) alkyl, alkenyl, alkynyl, carbocyclyl, heterocyclyl, etc.; R<sub>5</sub> = H, F, haloalkyl, (substituted) alkyl, alkenyl, alkynyl, heterocyclyl(alkyl), etc.; R<sub>13</sub> = H, F, alkyl, aminoalkyl, CF<sub>3</sub>, aminocarbonyl, etc.; R<sub>14</sub> = H, alkyl, aminoalkyl, F, CF<sub>3</sub>, aminocarbonyl, etc.; R<sub>13</sub>R<sub>14</sub> = O; R<sub>15</sub> = H, alkyl; R<sub>16</sub> = H, alkyl, PhCH<sub>2</sub>, alkylcarbonyl, alkylsulfonyl, alkoxy carbonyl], were prepared. Thus, 4-amidinobenzimidine monohydrochloride, styrene,

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1'-formyl-1-benzyloxycarbonyl-4-isobutylcarbamoylbiphenyl (preparation given) and indium triflate were heated together at 70° in MeCN for 12 h to give a product which was hydrogenolyzed in MeOH/HOAc over Pd/C to give 2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydroquinolin-2-yl)-4-isobutylcarbamoylbiphenyl-2-carboxylic acid. I inhibited Factor XIa with  $K_i \leq 15 \mu\text{M}$ .

AN 2004:780670 CAPLUS

DN 141:295874

TI Preparation of tetrahydroquinoline derivatives as inhibitors of serine protease enzymes of the coagulation cascade and/or contact activation system.

IN Quan, Mimi L.; Wang, Cailan; Zhou, Jinglan; Hangeland, Jon J.; Seiffert, Dietmar A.; Knabb, Robert M.

PA Bristol-Myers Squibb Company, USA

SO PCT Int. Appl., 150 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004080971	A1	20040923	WO 2004-US7216	20040310
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
	RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	US 2004235847	A1	20041125	US 2004-796396	20040309
PRAI	US 2003-453812P	P	20030311		
	US 2004-796396	A	20040309		
OS	MARPAT	141:295874			
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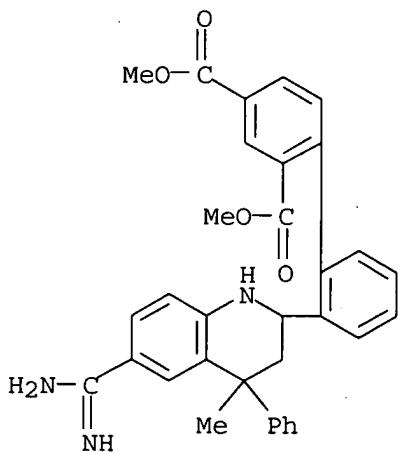
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of carbaminidoyltetrahydroquinoline derivs. as inhibitors of serine protease enzymes of the coagulation cascade and/or contact activation system)

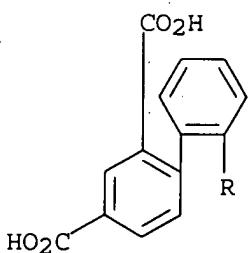
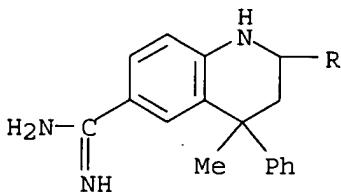
RN 762253-33-8 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 2'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-, dimethyl ester (9CI) (CA INDEX NAME)



RN 762253-34-9 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 2'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]- (9CI) (CA INDEX NAME)

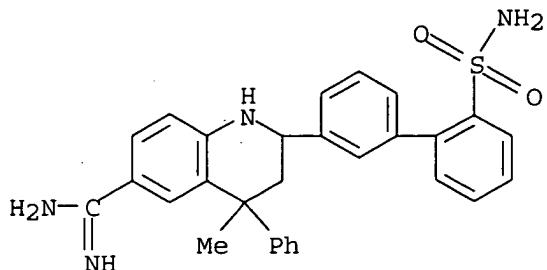


RN 762253-40-7 CAPLUS

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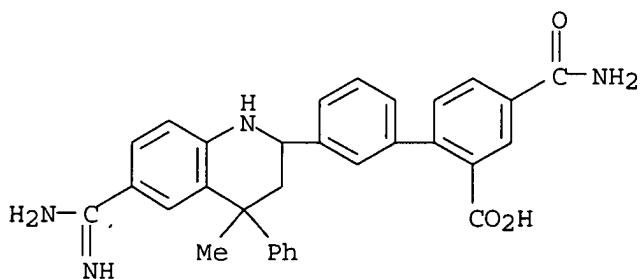
4/12/05

CN 6-Quinolinecarboximidamide, 2-[2'-(aminosulfonyl)[1,1'-biphenyl]-3-yl]-1,2,3,4-tetrahydro-4-methyl-4-phenyl- (9CI) (CA INDEX NAME)



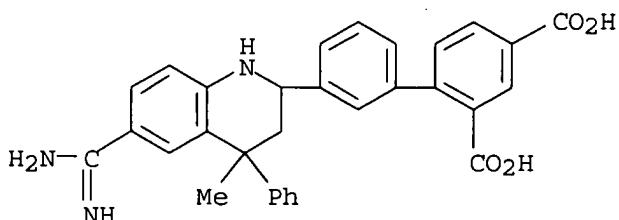
RN 762253-43-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)



RN 762253-45-2 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

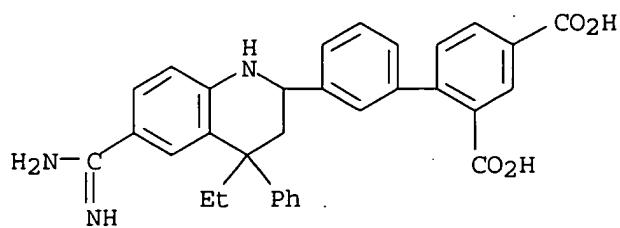


RN 762253-46-3 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 3'-(6-(aminoiminomethyl)-4-ethyl-1,2,3,4-tetrahydro-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

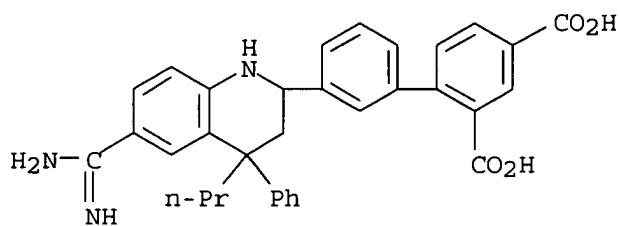
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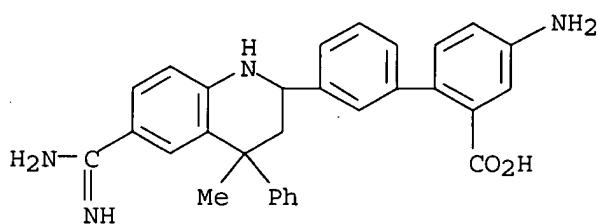
RN 762253-47-4 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-phenyl-4-propyl-2-quinoliny]- (9CI) (CA INDEX NAME)



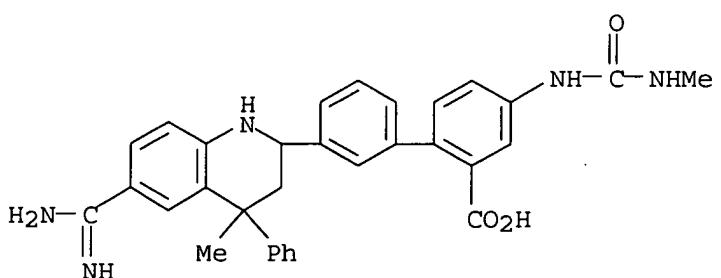
RN 762253-48-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-amino-3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinoliny]- (9CI) (CA INDEX NAME)



RN 762253-50-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinoliny]-4-[(methylamino)carbonyl]amino- (9CI) (CA INDEX NAME)

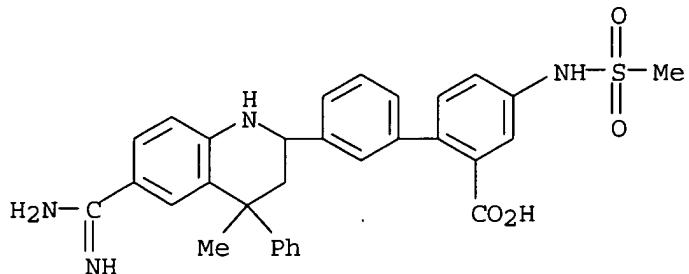


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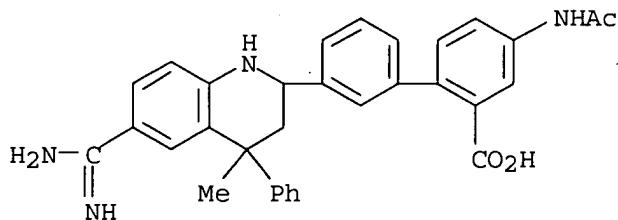
RN 762253-51-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4-[(methylsulfonyl)amino]-(9CI) (CA INDEX NAME)



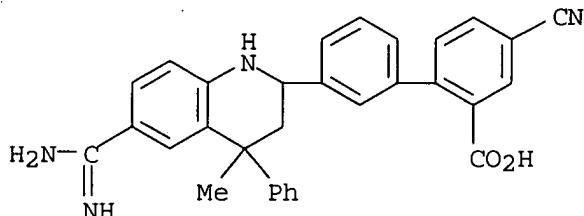
RN 762253-52-1 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(acetylamino)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-(9CI) (CA INDEX NAME)



RN 762253-53-2 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4-cyano-(9CI) (CA INDEX NAME)

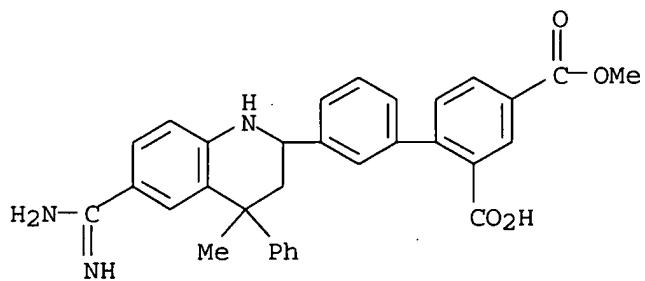


RN 762253-54-3 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, 4-methyl ester (9CI) (CA INDEX NAME)

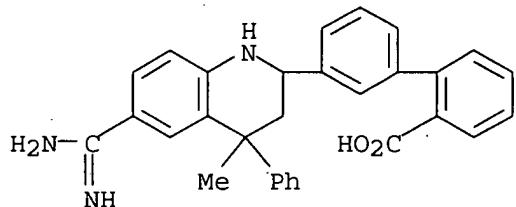
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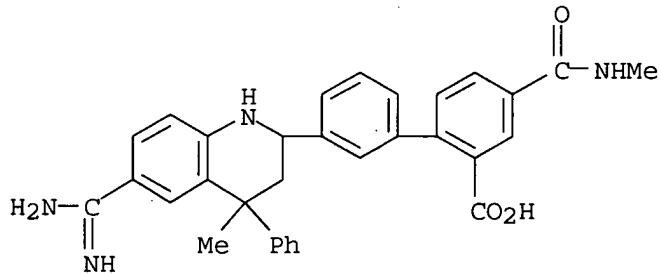
RN 762253-55-4 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]- (9CI) (CA INDEX NAME)



RN 762253-56-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-4-[(methylamino)carbonyl]- (9CI) (CA INDEX NAME)

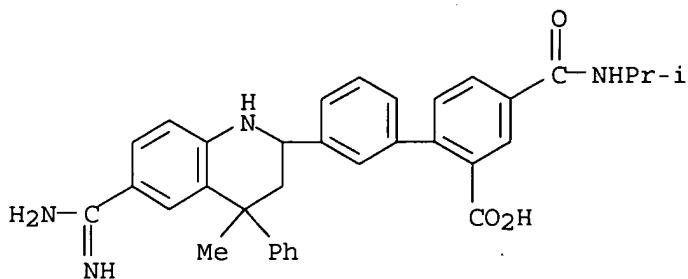


RN 762253-57-6 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-4-[(1-methylethyl)amino]carbonyl- (9CI) (CA INDEX NAME)

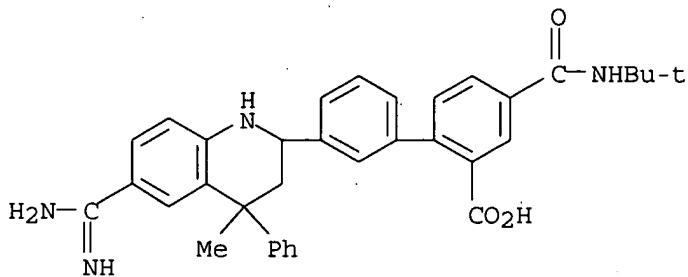
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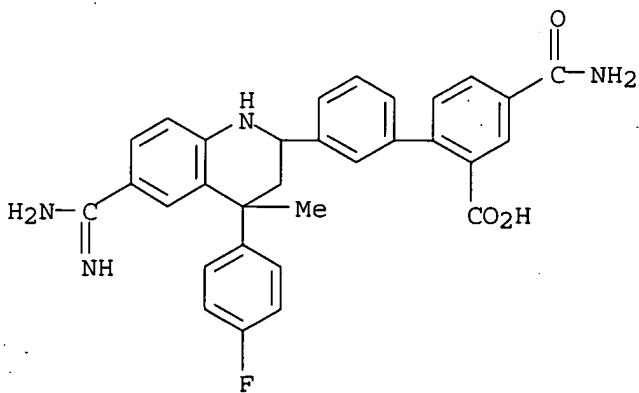
RN 762253-58-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-4-[(1,1-dimethylethyl)amino]carbonyl- (9CI) (CA INDEX NAME)



RN 762253-59-8 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-4-(4-fluorophenyl)-1,2,3,4-tetrahydro-4-methyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

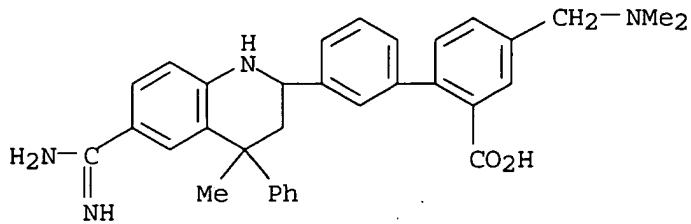


RN 762253-60-1 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4-[(dimethylamino)methyl]- (9CI) (CA INDEX NAME)

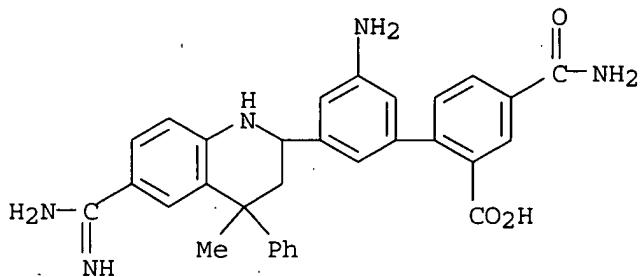
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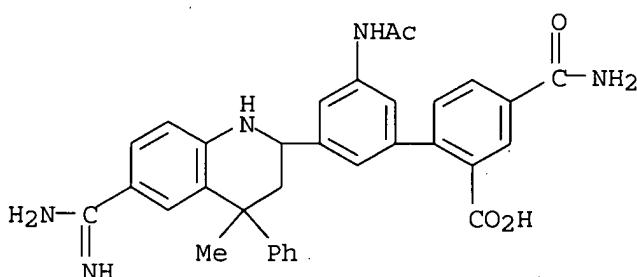
RN 762253-62-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-amino-4-(aminocarbonyl)-5'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]- (9CI) (CA INDEX NAME)



RN 762253-64-5 CAPLUS

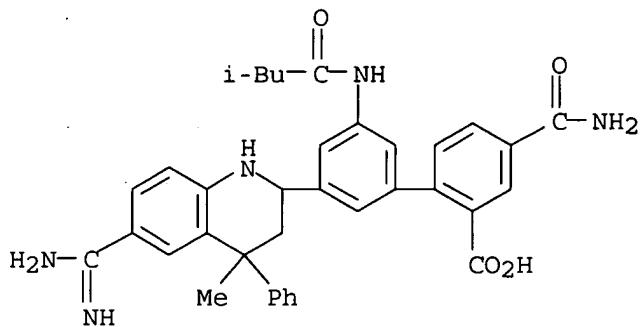
CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(acetylamino)-4-(aminocarbonyl)-5'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]- (9CI) (CA INDEX NAME)



RN 762253-65-6 CAPLUS

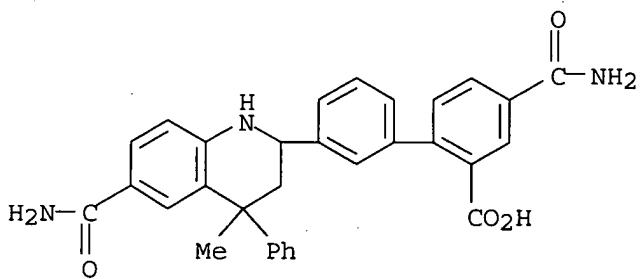
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-5'-[(3-methyl-1-oxobutyl)amino]- (9CI) (CA INDEX NAME)

4/12/05



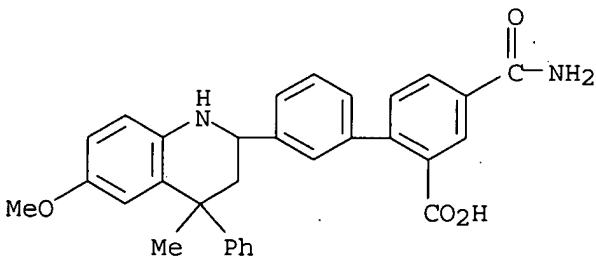
RN 762253-66-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminocarbonyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)



RN 762253-67-8 CAPLUS

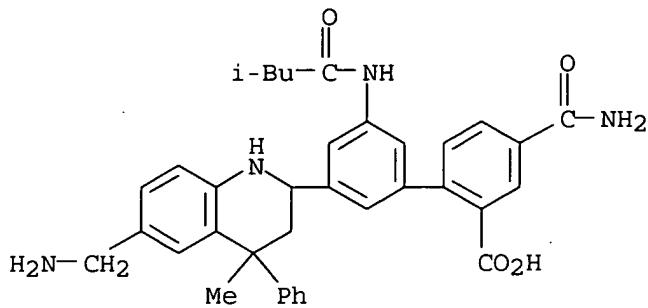
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(1,2,3,4-tetrahydro-6-methoxy-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)



RN 762253-68-9 CAPLUS

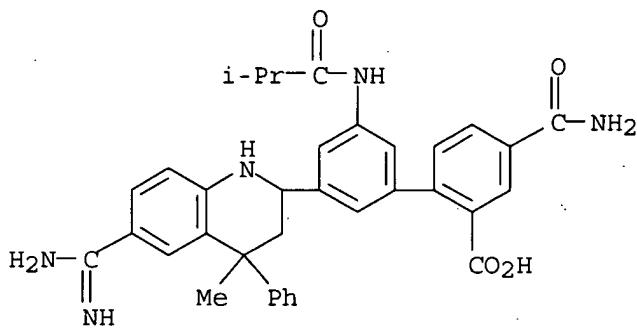
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(3-methyl-1-oxobutylamino)- (9CI) (CA INDEX NAME)

4/12/05



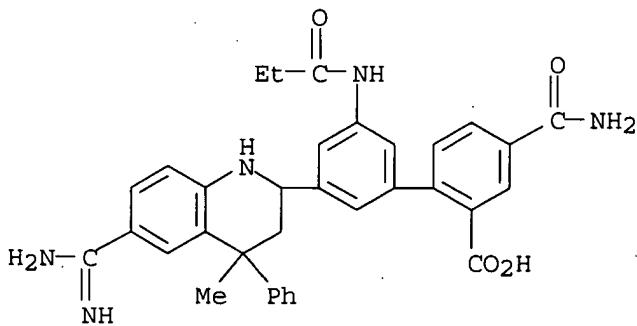
RN 762253-69-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(2-methyl-1-oxopropyl)amino] - (9CI) (CA INDEX NAME)



RN 762253-70-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(1-oxopropyl)amino] - (9CI) (CA INDEX NAME)

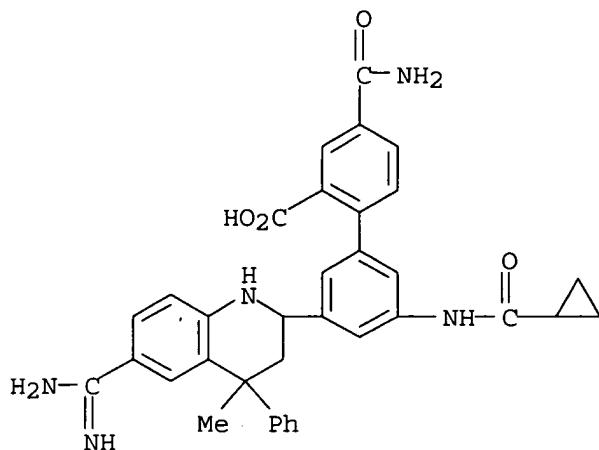


RN 762253-71-4 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(cyclopropylcarbonyl)amino] - (9CI) (CA INDEX NAME)

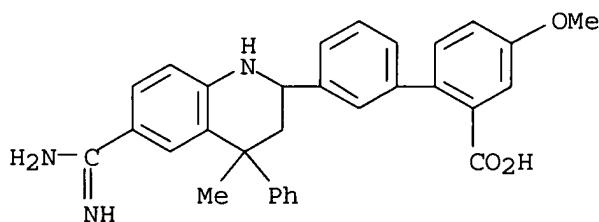
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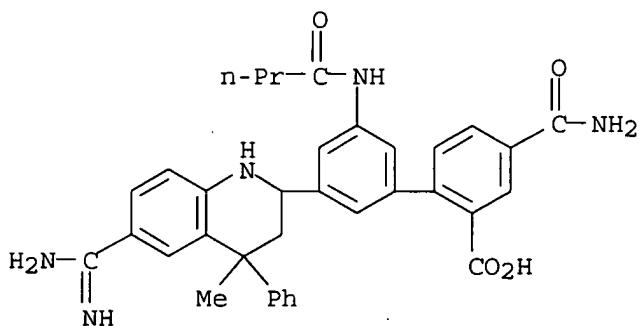
RN 762253-72-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-4-methoxy- (9CI) (CA INDEX NAME)



RN 762253-73-6 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5-[(1-oxobutyl)amino]- (9CI) (CA INDEX NAME)

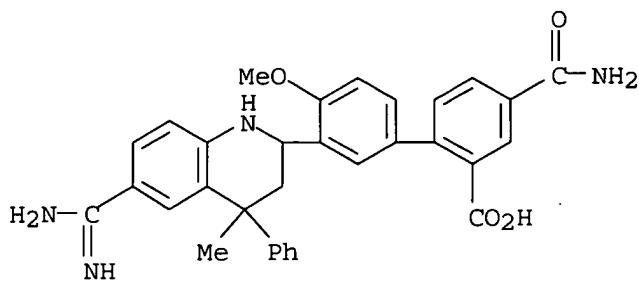


RN 762253-74-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4'-methoxy- (9CI) (CA INDEX NAME)

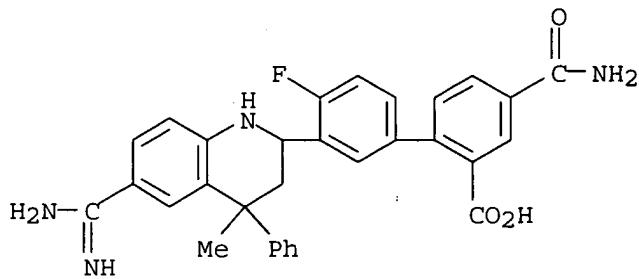
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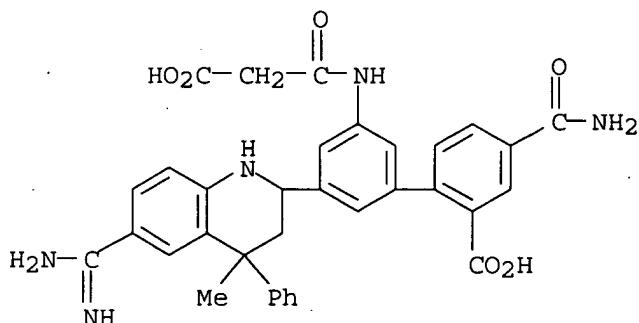
RN 762253-75-8 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4'-fluoro- (9CI) (CA INDEX NAME)



RN 762253-76-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[(carboxyacetyl)amino]- (9CI) (CA INDEX NAME)

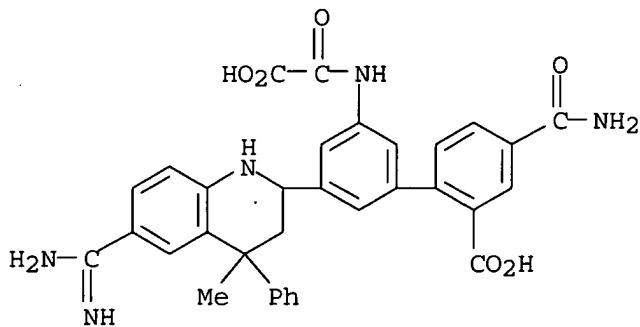


RN 762253-77-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[(carboxycarbonyl)amino]- (9CI) (CA INDEX NAME)

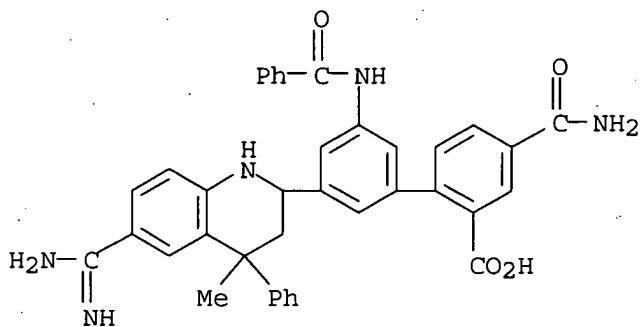
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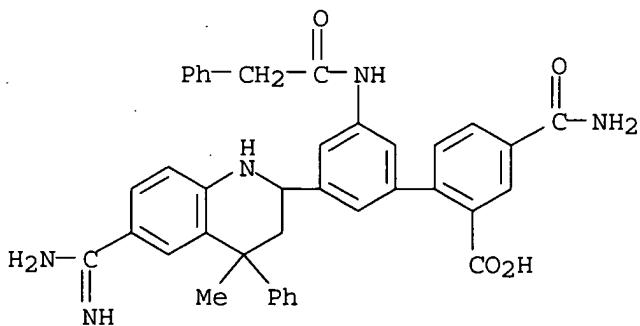
RN 762253-78-1 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(benzoylamino)- (9CI) (CA INDEX NAME)



RN 762253-79-2 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-([(phenylacetyl)amino]- (9CI) (CA INDEX NAME)

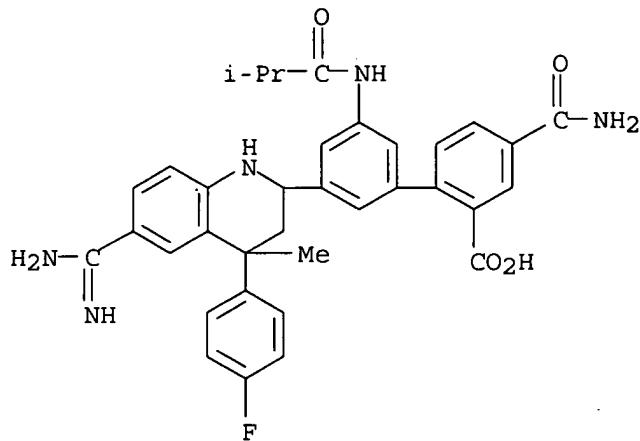


RN 762253-80-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-4-(4-fluorophenyl)-1,2,3,4-tetrahydro-4-methyl-2-quinolinyl)-5'-([(2-methyl-1-oxopropyl)amino]- (9CI) (CA INDEX NAME)

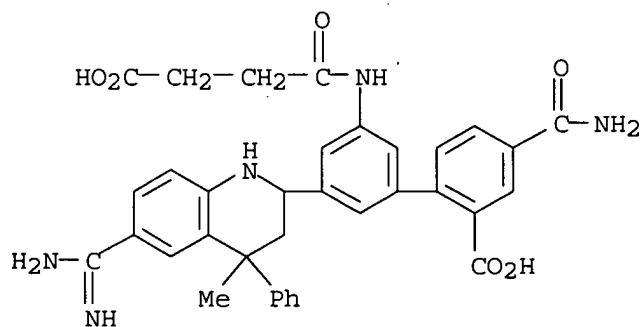
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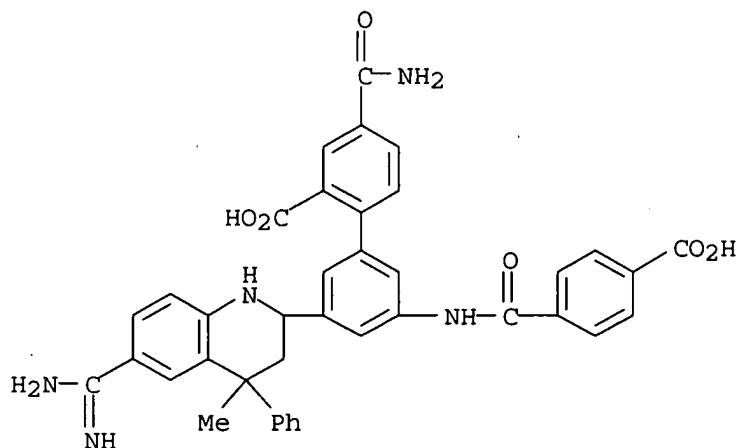
RN 762253-81-6 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(3-carboxy-1-oxopropyl)amino] - (9CI) (CA INDEX NAME)



RN 762253-82-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(4-carboxybenzoyl)amino] - (9CI) (CA INDEX NAME)

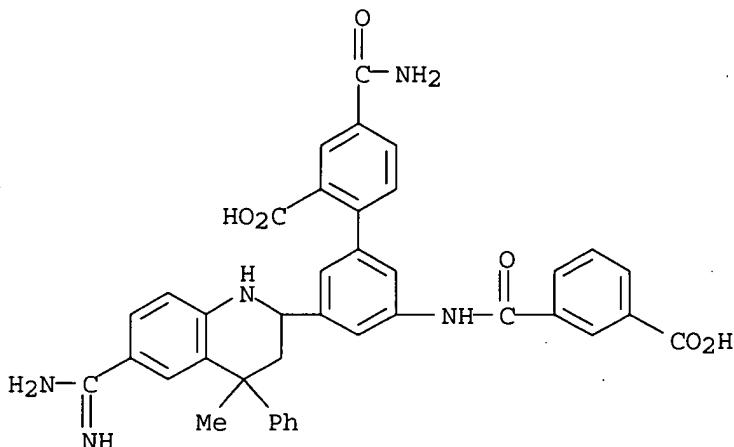


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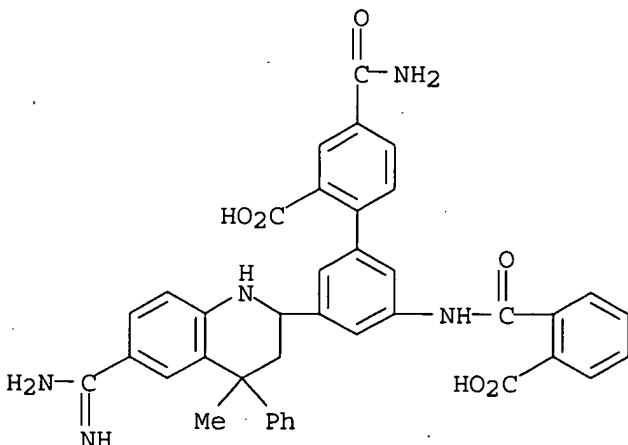
RN 762253-83-8 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(3-carboxybenzoyl)amino] - (9CI) (CA INDEX NAME)



RN 762253-84-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(2-carboxybenzoyl)amino] - (9CI) (CA INDEX NAME)

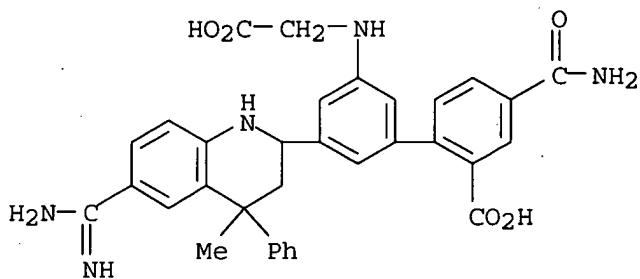


RN 762253-85-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(carboxymethyl)amino] - (9CI) (CA INDEX NAME)

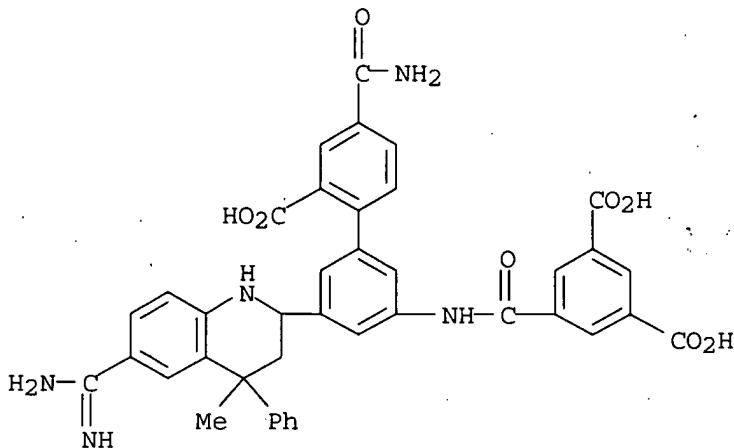
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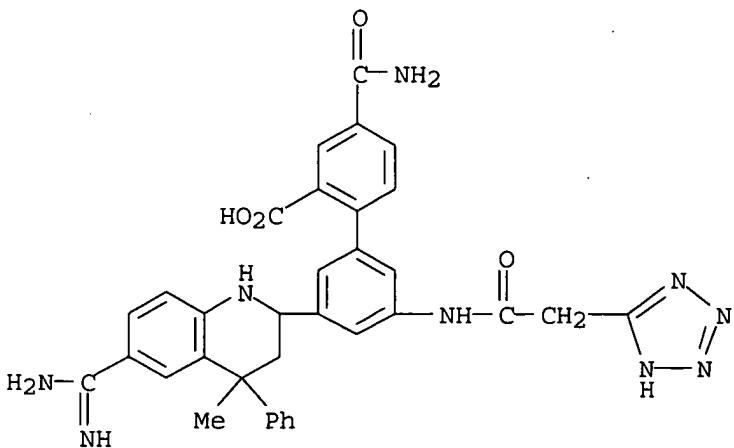
RN 762253-86-1 CAPLUS

CN 1,3-Benzenedicarboxylic acid, 5-[[[4'-(aminocarbonyl)-5-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-2'-carboxy[1,1'-biphenyl]-3-yl]amino]carbonyl]- (9CI) (CA INDEX NAME)



RN 762253-87-2 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-5'-[(1H-tetrazol-5-ylacetyl)amino]- (9CI) (CA INDEX NAME)

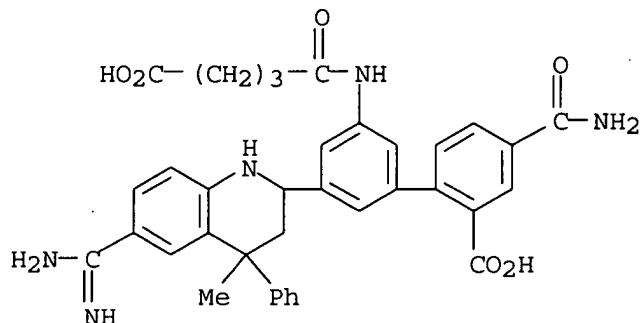


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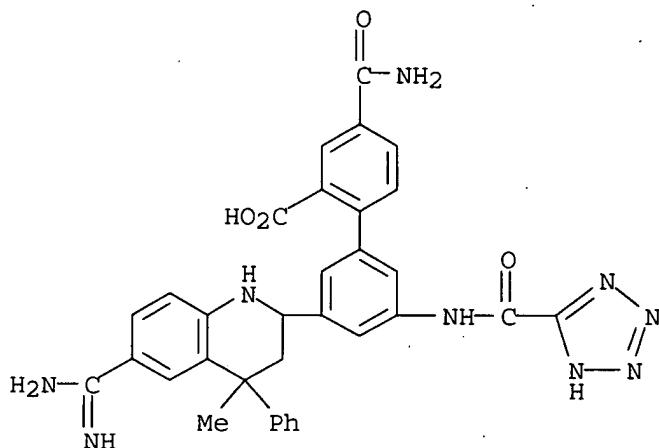
RN 762253-88-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-([(4-carboxy-1-oxobutyl)amino]- (9CI) (CA INDEX NAME)



RN 762253-89-4 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-([(1H-tetrazol-5-ylcarbonyl)amino]- (9CI) (CA INDEX NAME)

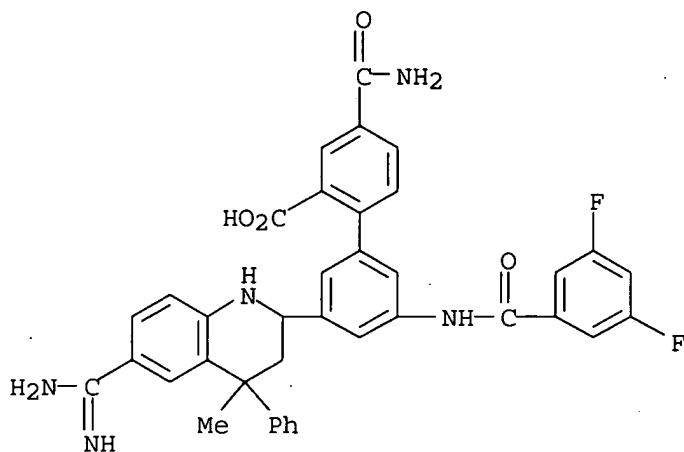


RN 762253-91-8 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-([(3,5-difluorobenzoyl)amino]- (9CI) (CA INDEX NAME)

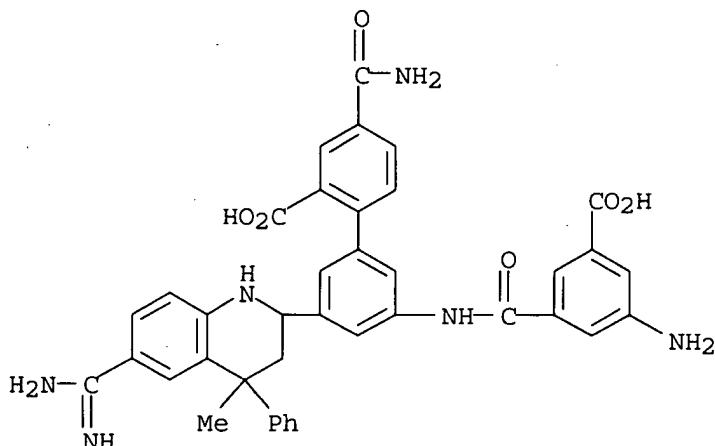
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RN 762253-92-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(3-amino-5-carboxybenzoyl)amino]-5'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

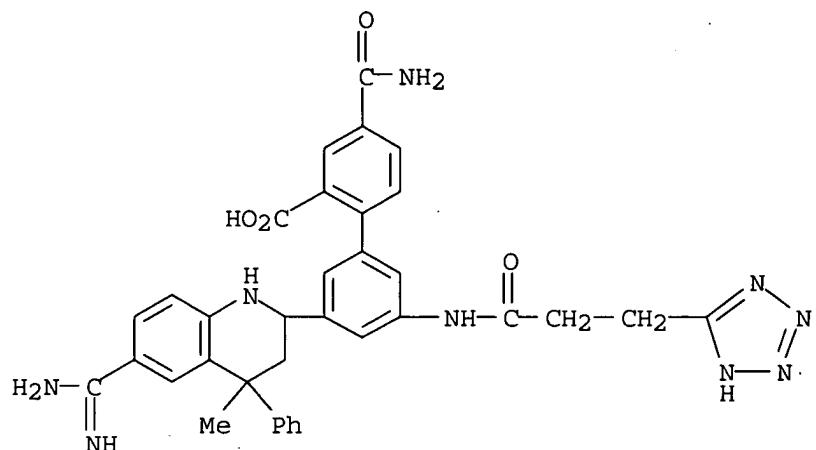


RN 762253-93-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(1-oxo-3-(1H-tetrazol-5-yl)propylamino)- (9CI) (CA INDEX NAME)

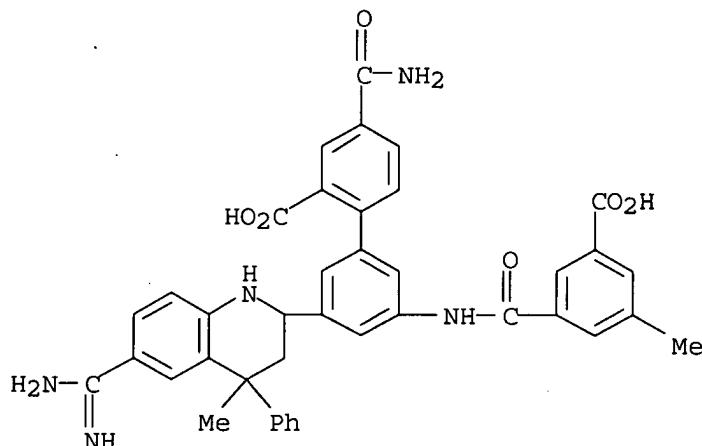
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RN 762253-94-1 CAPLUS

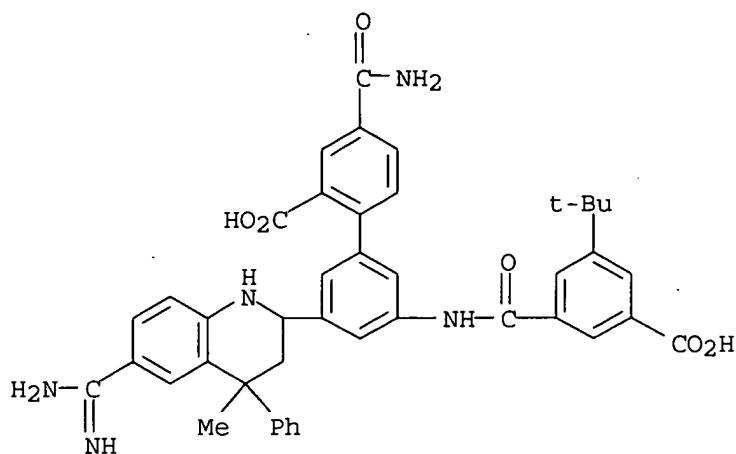
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[(3-carboxy-5-methylbenzoyl)amino]- (9CI) (CA INDEX NAME)



RN 762253-95-2 CAPLUS

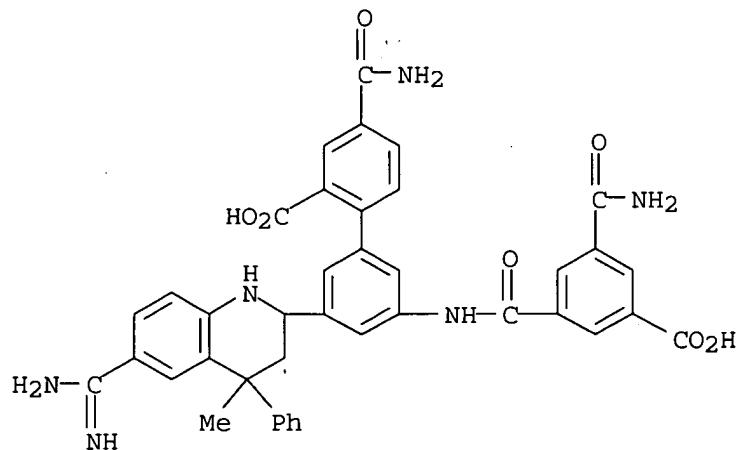
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[(3-carboxy-5-(1,1-dimethylethyl)benzoyl)amino]- (9CI) (CA INDEX NAME)

4/12/05



RN 762253-96-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(3-(aminocarbonyl)-5-carboxybenzoyl)amino]-5'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

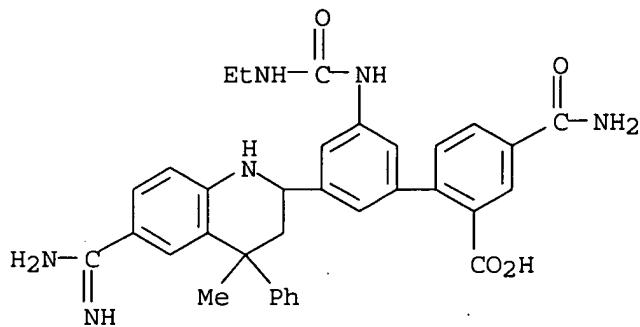


RN 762253-97-4 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(ethylamino)carbonyl)amino]-5'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)

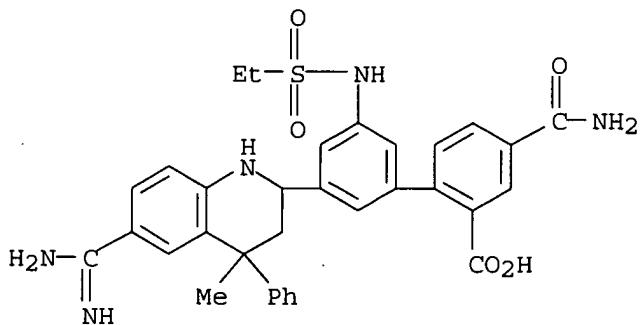
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RN 762253-98-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(ethylsulfonyl)amino]-(9CI) (CA INDEX NAME)



RN 762254-03-5 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 2'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, dimethyl ester, trifluoroacetate (9CI) (CA INDEX NAME)

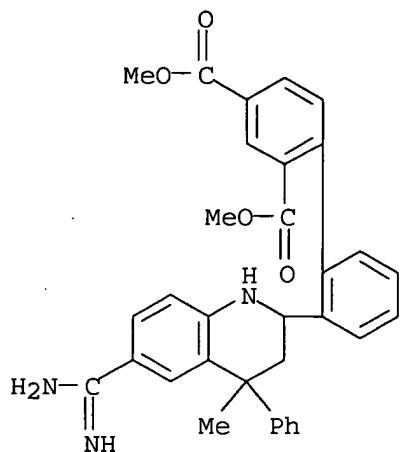
CM 1

CRN 762253-33-8

CMF C33 H31 N3 O4

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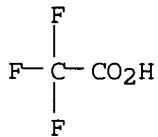
4/12/05



CM 2

CRN 76-05-1

CMF C2 H F3 O2



RN 762254-04-6 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 2'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-, trifluoroacetate (9CI) (CA INDEX NAME)

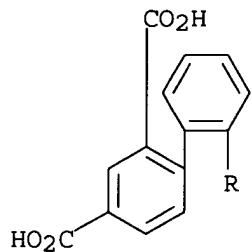
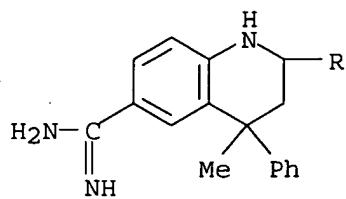
CM 1

CRN 762253-34-9

CMF C31 H27 N3 O4

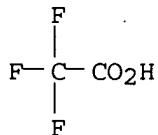
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CM 2

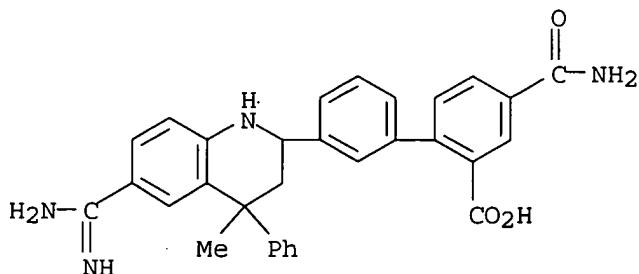
CRN 76-05-1  
CMF C2 H F3 O2



RN 762254-08-0 CAPLUS  
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762253-43-0  
CMF C31 H28 N4 O3

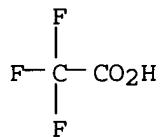


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CM 2

CRN 76-05-1  
CMF C2 H F3 O2

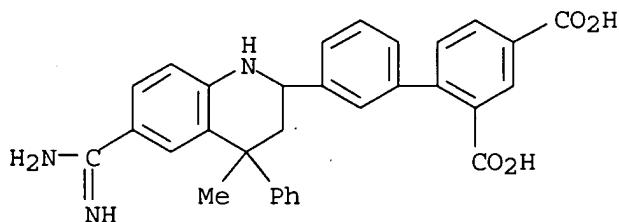


RN 762254-10-4 CAPLUS

CN [1,1'-Biphenyl]-2,4-dicarboxylic acid, 3'-[6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-, trifluoroacetate (9CI) (CA INDEX NAME)

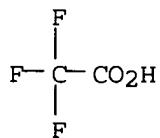
CM 1

CRN 762253-45-2  
CMF C31 H27 N3 O4



CM 2

CRN 76-05-1  
CMF C2 H F3 O2



RN 762254-11-5 CAPLUS

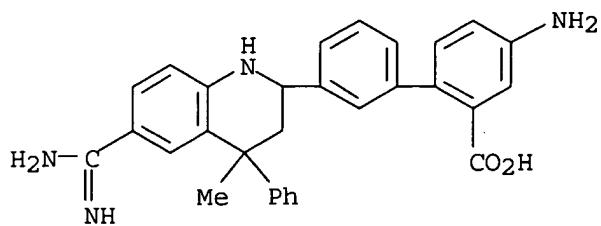
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-amino-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762253-48-5  
CMF C30 H28 N4 O2

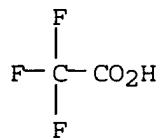
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4/12/05



CM 2

CRN 76-05-1  
CMF C2 H F3 O2

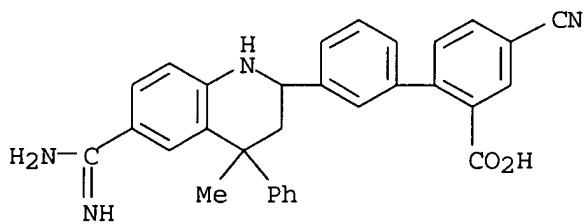


RN 762254-12-6 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4-cyano-, trifluoroacetate (9CI) (CA INDEX NAME)

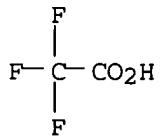
CM 1

CRN 762253-53-2  
CMF C31 H26 N4 O2



CM 2

CRN 76-05-1  
CMF C2 H F3 O2



10796396

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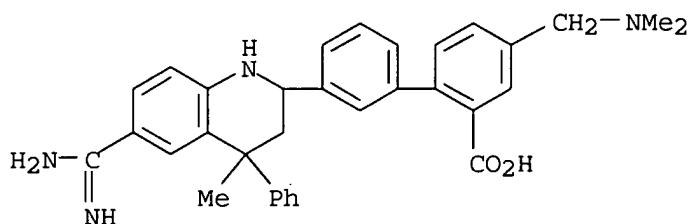
RN 762254-13-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-4-[(dimethylamino)methyl]-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762253-60-1

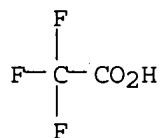
CMF C33 H34 N4 O2



CM 2

CRN 76-05-1

CMF C2 H F3 O2



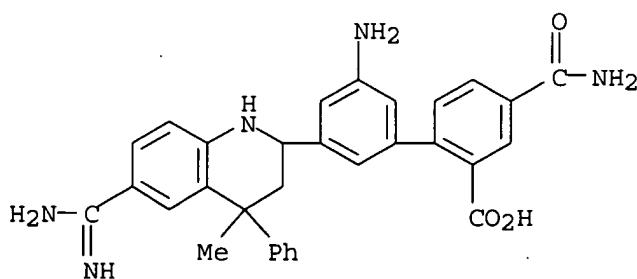
RN 762254-15-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-amino-4-(aminocarbonyl)-5'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762253-62-3

CMF C31 H29 N5 O3

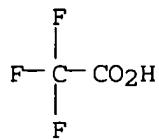


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CM 2

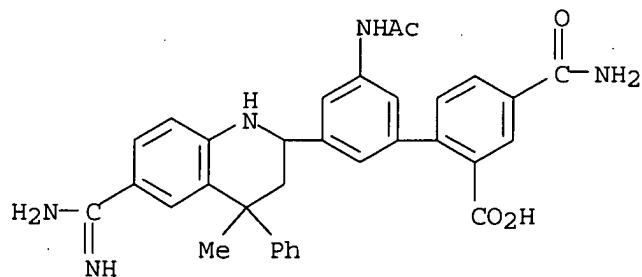
CRN 76-05-1  
CMF C2 H F3 O2



RN 762254-16-0 CAPLUS  
CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(acetylamino)-4-(aminocarbonyl)-5'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

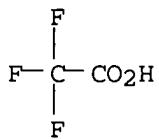
CM 1

CRN 762253-64-5  
CMF C33 H31 N5 O4



CM 2

CRN 76-05-1  
CMF C2 H F3 O2



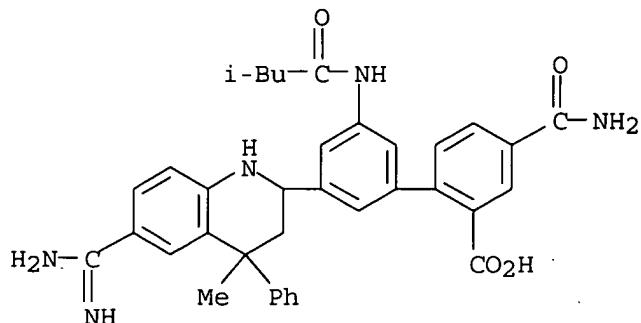
RN 762254-17-1 CAPLUS  
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-(3-methyl-1-oxobutyl)amino-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

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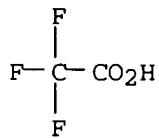
4/12/05

CRN 762253-65-6  
CMF C36 H37 N5 O4



CM 2

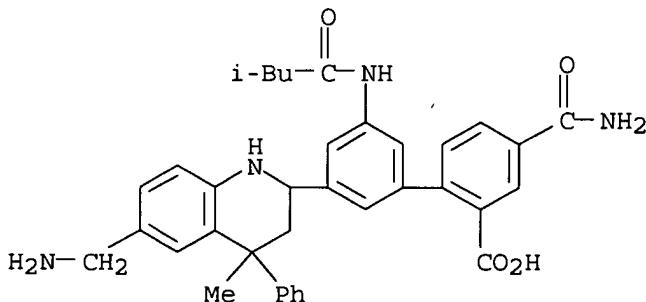
CRN 76-05-1  
CMF C2 H F3 O2



RN 762254-19-3 CAPLUS  
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-[6-(aminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl]-5'-(3-methyl-1-oxobutyl)amino]-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762253-68-9  
CMF C36 H38 N4 O4

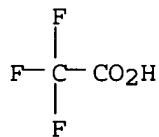


CM 2

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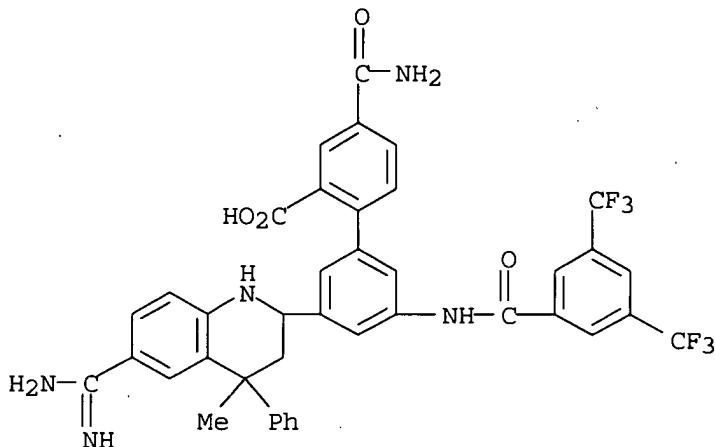
4/12/05

CRN 76-05-1  
CMF C2 H F3 O2



RN 762254-69-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-(aminoiminomethyl)-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[[3,5-bis(trifluoromethyl)benzoyl]amino]-(9CI) (CA INDEX NAME)



IT 762254-30-8P 762254-31-9P 762254-32-0P  
762254-34-2P 762254-35-3P 762254-37-5P  
762254-39-7P 762254-47-7P 762254-50-2P  
762254-58-0P 762254-60-4P 762254-61-5P  
762254-62-6P 762254-63-7P 762254-65-9P  
762254-67-1P

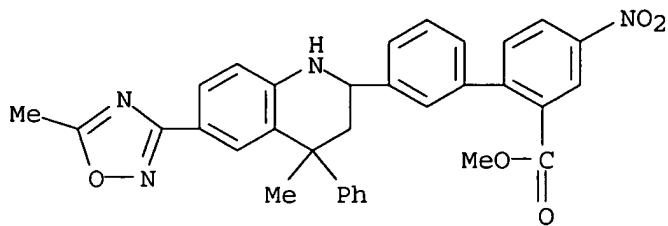
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of carbaminidoyltetrahydroquinoline derivs. as inhibitors of serine protease enzymes of the coagulation cascade and/or contact activation system)

RN 762254-30-8 CAPLUS

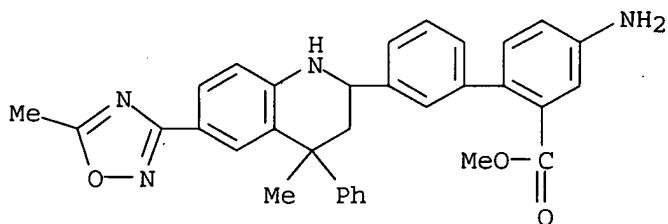
CN [1,1'-Biphenyl]-2-carboxylic acid, 4-nitro-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)

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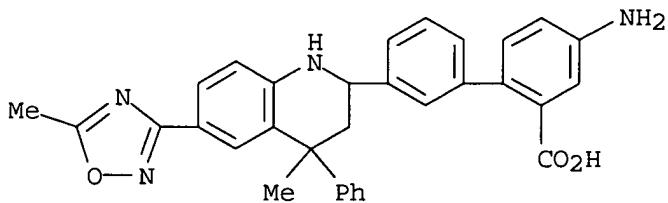
RN 762254-31-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-amino-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)



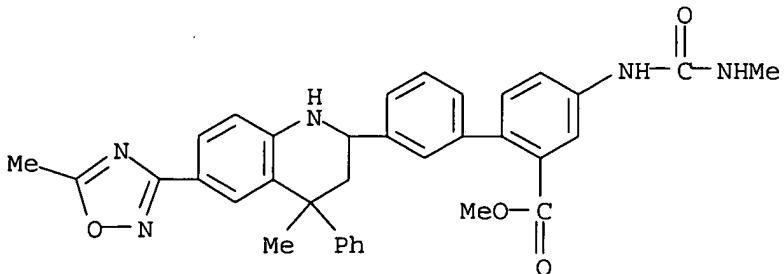
RN 762254-32-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-amino-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)- (9CI) (CA INDEX NAME)



RN 762254-34-2 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-[(methylamino)carbonyl]amino-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)

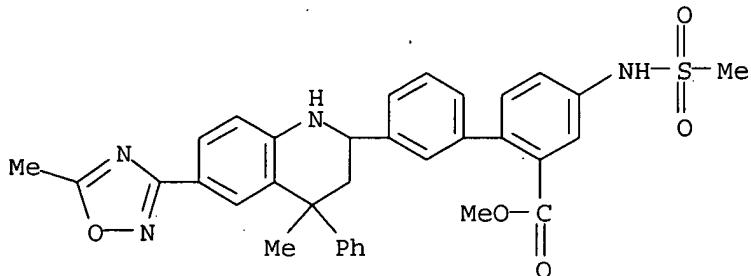


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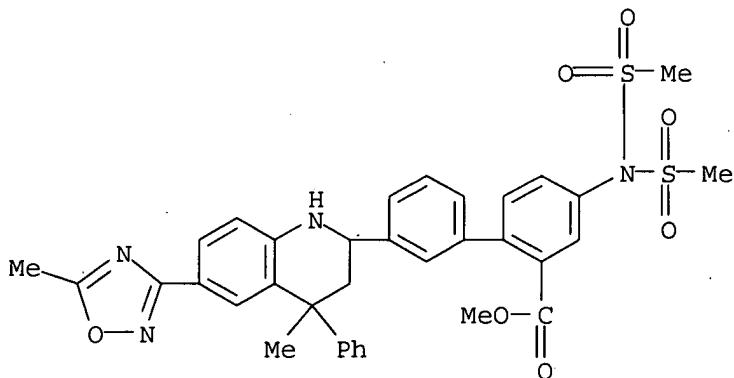
RN 762254-35-3 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-[(methylsulfonyl)amino]-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)



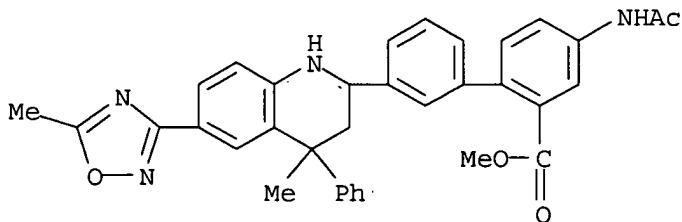
RN 762254-37-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-[bis(methylsulfonyl)amino]-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)



RN 762254-39-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(acetylamino)-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, methyl ester (9CI) (CA INDEX NAME)



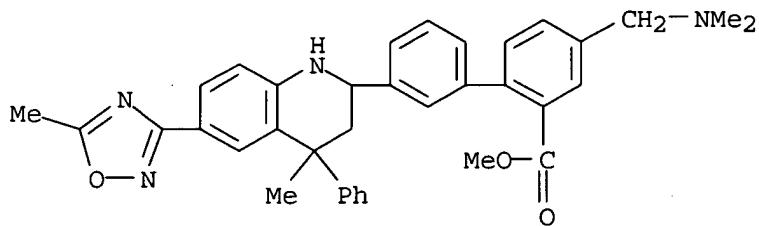
RN 762254-47-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-[(dimethylamino)methyl]-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-

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quinolinyl]-, methyl ester (9CI) (CA INDEX NAME)



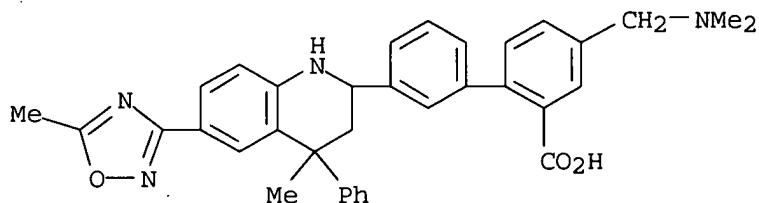
RN 762254-50-2 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-[(dimethylamino)methyl]-3'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 762254-49-9

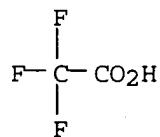
CMF C35 H34 N4 O3



CM 2

CRN 76-05-1

CMF C2 H F3 O2

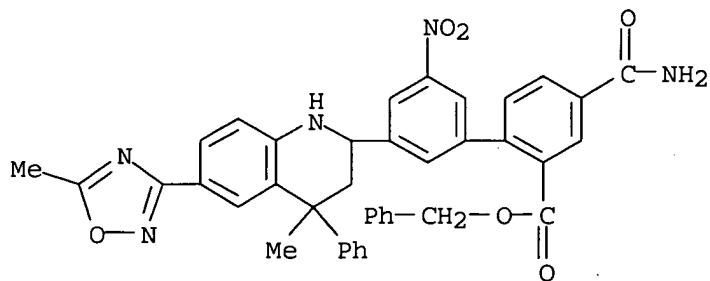


RN 762254-58-0 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-nitro-5'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, phenylmethyl ester (9CI) (CA INDEX NAME)

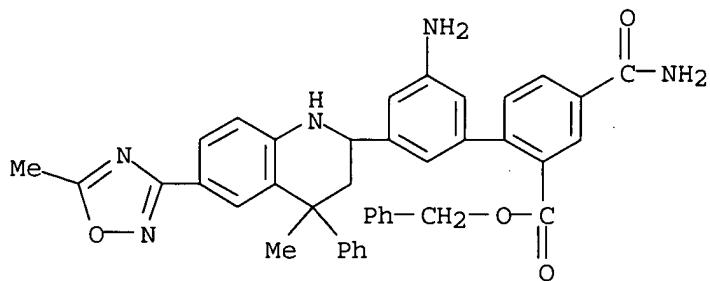
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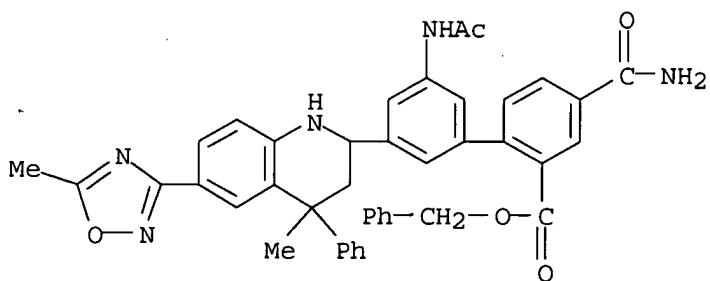
RN 762254-60-4 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-amino-4-(aminocarbonyl)-5'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, phenylmethyl ester (9CI) (CA INDEX NAME)



RN 762254-61-5 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-(acetylamino)-4-(aminocarbonyl)-5'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, phenylmethyl ester (9CI) (CA INDEX NAME)

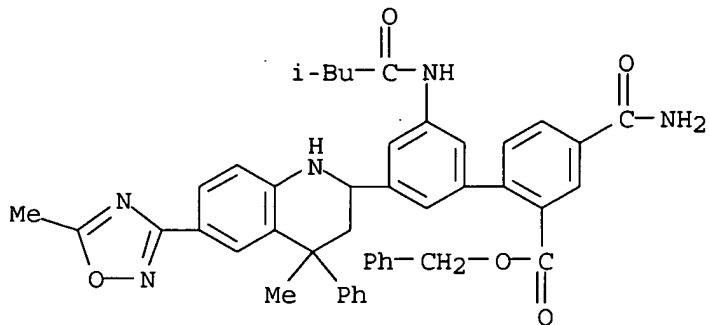


RN 762254-62-6 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(3-methyl-1-oxobutyl)amino-5'-(1,2,3,4-tetrahydro-4-methyl-6-(5-methyl-1,2,4-oxadiazol-3-yl)-4-phenyl-2-quinolinyl)-, phenylmethyl ester (9CI) (CA INDEX NAME)

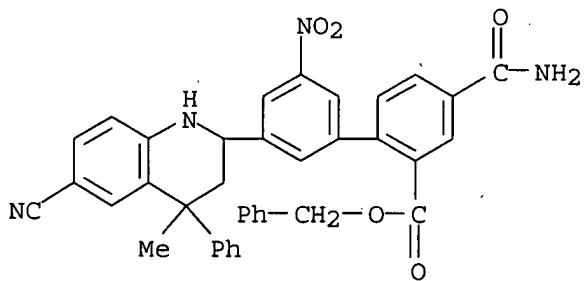
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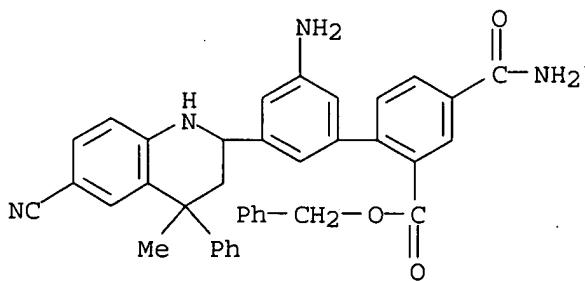
RN 762254-63-7 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-cyano-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-nitro-, phenylmethyl ester (9CI) (CA INDEX NAME)



RN 762254-65-9 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 3'-amino-4-(aminocarbonyl)-5'-(6-cyano-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-, phenylmethyl ester (9CI) (CA INDEX NAME)

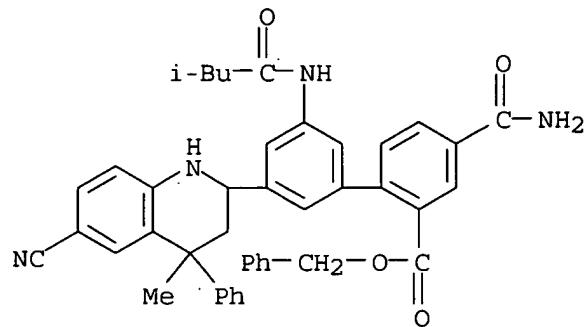


RN 762254-67-1 CAPLUS

CN [1,1'-Biphenyl]-2-carboxylic acid, 4-(aminocarbonyl)-3'-(6-cyano-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-quinolinyl)-5'-[(3-methyl-1-oxobutyl)amino]-, phenylmethyl ester (9CI) (CA INDEX NAME)

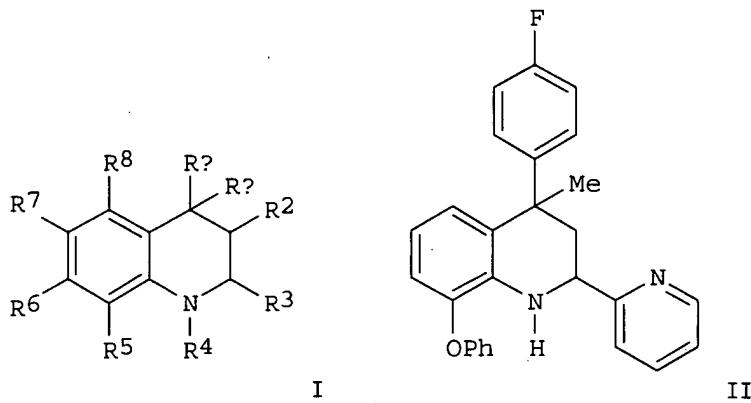
10796396

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RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
GI



AB Title compds. I [Rb and R2 together =  $(CH_2)_n$ ,  $CH=CHCH_2$ ,  $CH_2CH=CH$ , etc.; n = 3-10; Ra = H; R3 = H, alkyl, cycloalkyl, etc.; R4 = R4a, ZR4a; R4a = H, alkyl, alkenyl, etc.; Z = alkyl, alkenyl, alkynyl, etc.; R5, R6, R7, R8 = H, halo, CN, etc.] and their pharmaceutically acceptable salts were prepared In sodium channel [<sup>3</sup>H]batrachotoxin (BTX) displacement assays, 261-examples of compds. I exhibited 00.0-91.7% binding, e.g., the affinity of tetrahydroquinoline was 91.7%.

AN 2004:198148 CAPLUS

DN 140:253456

TI Preparation of 1,2,3,4-tetrahydro-4-phenylquinolines and related compounds as sodium channel ligands for the treatment of pain

IN Hennies, Hagen-Heinrich; Maul, Corinna; Przewosny, Michael; Sundermann, Bernd

PA Gruenenthal G.m.b.H., Germany

SO Ger. Offen., 55 pp.

CODEN: GWXXBX

DT Patent

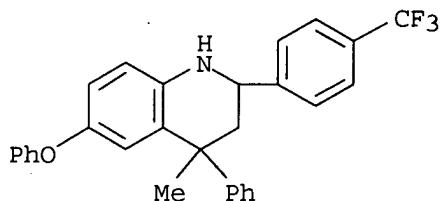
LA German

FAN.CNT 1

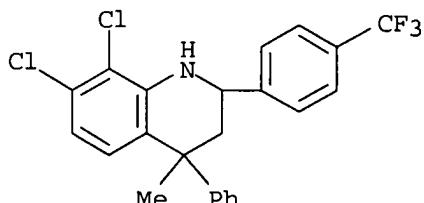
10796396

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	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	DE 10236910 WO 2004022542 WO 2004022542	A1 A2 A3	20040311 20040318 20040603	DE 2002-10236910 WO 2003-EP8889	20020812 20030811
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
PRAI	DE 2002-10236910	A	20020812		
OS	MARPAT 140:253456				
IT	669778-57-8 669778-87-4 669778-95-4 669779-11-7 669779-28-6 669779-35-5 669779-37-7 669779-45-7 669779-86-6 669779-98-0 669780-06-7				
	RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses) (preparation of tetrahydroquinolines and related compds. as sodium channel ligands for the treatment of pain)				
RN	669778-57-8 CAPLUS				
CN	Quinoline, 1,2,3,4-tetrahydro-4-methyl-6-phenoxy-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)				



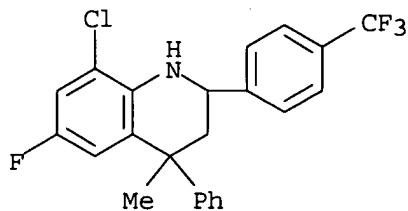
RN 669778-87-4 CAPLUS  
CN Quinoline, 7,8-dichloro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 669778-95-4 CAPLUS  
CN Quinoline, 8-chloro-6-fluoro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-[4-(trifluoromethyl)phenyl].- (9CI) (CA INDEX NAME)

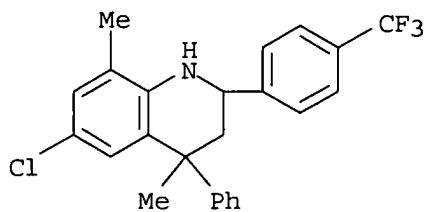
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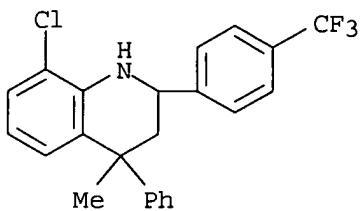
RN 669779-11-7 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4,8-dimethyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



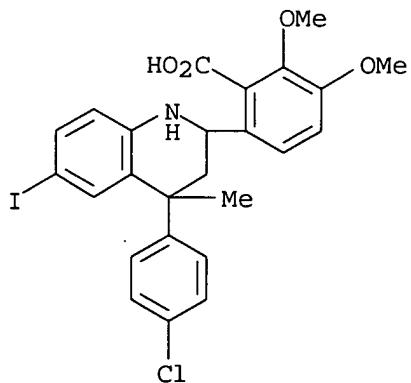
RN 669779-28-6 CAPLUS

CN Quinoline, 8-chloro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 669779-35-5 CAPLUS

CN Benzoic acid, 6-[4-(4-chlorophenyl)-1,2,3,4-tetrahydro-6-iodo-4-methyl-2-quinolinyl]-2,3-dimethoxy- (9CI) (CA INDEX NAME)

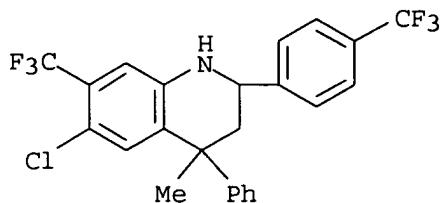


RN 669779-37-7 CAPLUS

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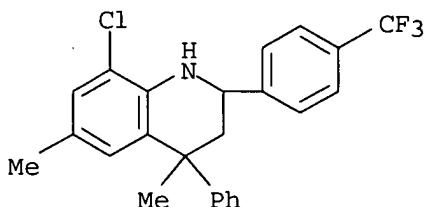
4/12/05

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-7-(trifluoromethyl)-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



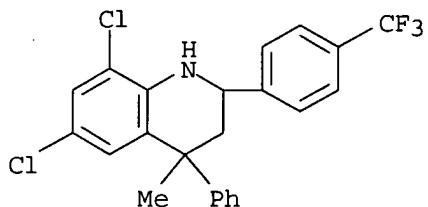
RN 669779-45-7 CAPLUS

CN Quinoline, 8-chloro-1,2,3,4-tetrahydro-4,6-dimethyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 669779-86-6 CAPLUS

CN Quinoline, 6,8-dichloro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)

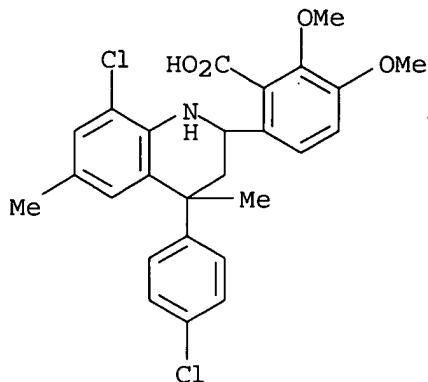


RN 669779-98-0 CAPLUS

CN Benzoic acid, 6-[8-chloro-4-(4-chlorophenyl)-1,2,3,4-tetrahydro-4,6-dimethyl-2-quinolinyl]-2,3-dimethoxy- (9CI) (CA INDEX NAME)

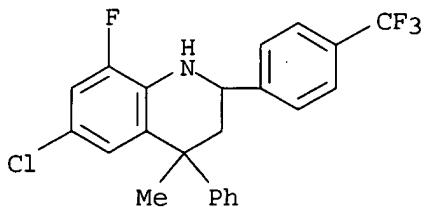
10796396

4/12/05



RN 669780-06-7 CAPLUS

CN Quinoline, 6-chloro-8-fluoro-1,2,3,4-tetrahydro-4-methyl-4-phenyl-2-[4-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



RE.CNT 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN

AB Cycloaddn. of N-arylimines with  $\alpha$ -methylstyrenes or 2,3-dihydrofuran was efficiently catalyzed by tris(4-bromophenyl)aminium hexachloroantimonate ( $\text{Ar}_3\text{N}^+\text{SbCl}_6^-$ ) producing tetrahydroquinoline derivs. in excellent yields. The reaction was controlled sensitively by the oxidation potentials of the imine and the dienophile.

AN 2003:732818 CAPLUS

DN 140:42014

TI Cation radical imino Diels-Alder reaction: A new approach for the synthesis of tetrahydroquinolines

AU Jia, Xiaodong; Lin, Hechun; Huo, Congde; Zhang, Wei; Lue, Jianming; Yang, Li; Zhao, Guangyu; Liu, Zhong-Li

CS National Laboratory of Applied Organic Chemistry, Lanzhou University, Lanzhou, 730000, Peop. Rep. China

SO Synlett (2003), (11), 1707-1709

CODEN: SYNLES; ISSN: 0936-5214

PB Georg Thieme Verlag

DT Journal

LA English

OS CASREACT 140:42014

IT 506407-52-9P 506407-53-0P 506407-54-1P

506407-64-3P 506407-65-4P 506407-66-5P

635303-04-7P 635303-05-8P 635303-06-9P

635303-07-0P 727653-41-0P 727653-65-8P

727653-68-1P 727653-73-8P

RL: SPN (Synthetic preparation); PREP (Preparation)

10796396

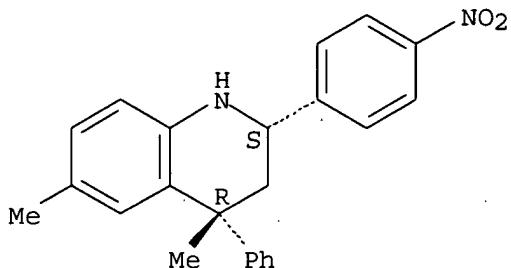
4/12/05

(preparation of tetrahydroquinolines via cation radical imino Diels-Alder reaction of N-arylimines with dienophiles catalyzed by tris(bromophenyl)aminium hexachloroantimonate)

RN 506407-52-9 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

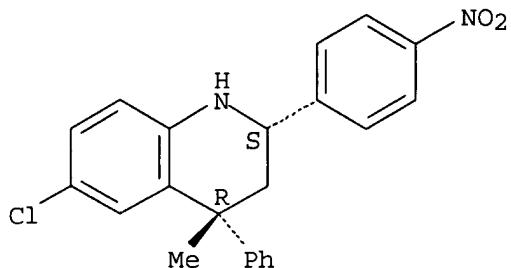
Relative stereochemistry.



RN 506407-53-0 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

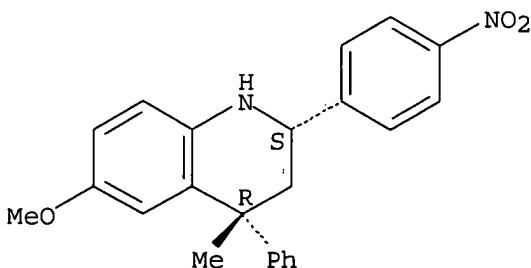
Relative stereochemistry.



RN 506407-54-1 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



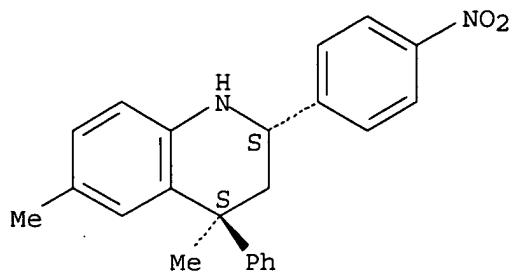
RN 506407-64-3 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

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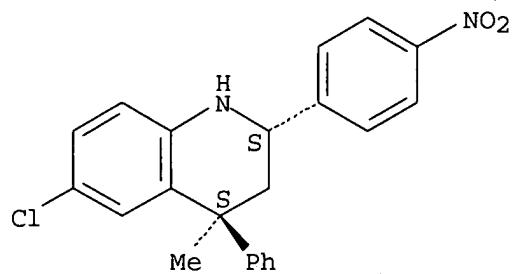
Relative stereochemistry.



RN 506407-65-4 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

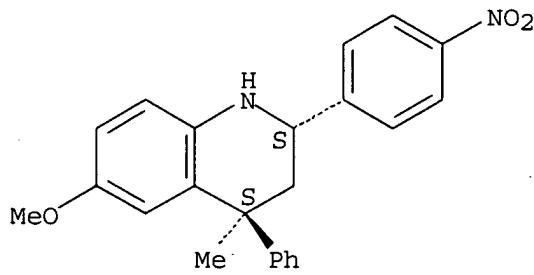
Relative stereochemistry.



RN 506407-66-5 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



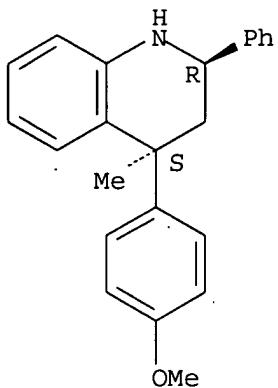
RN 635303-04-7 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4-(4-methoxyphenyl)-4-methyl-2-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

10796396

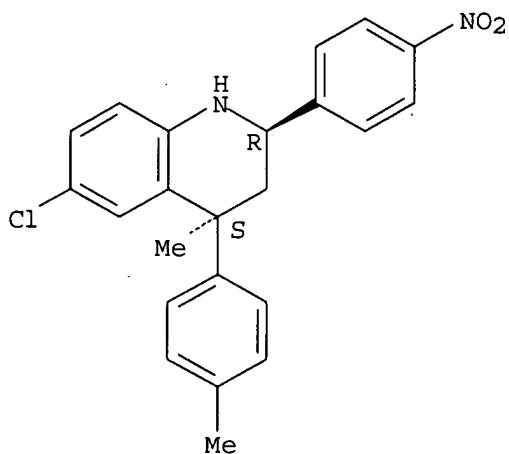
4/12/05



RN 635303-05-8 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-4-(4-methylphenyl)-2-(4-nitrophenyl)-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

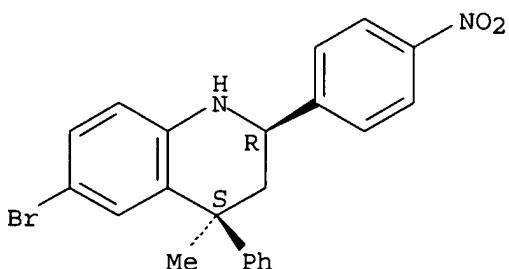
Relative stereochemistry.



RN 635303-06-9 CAPLUS

CN Quinoline, 6-bromo-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



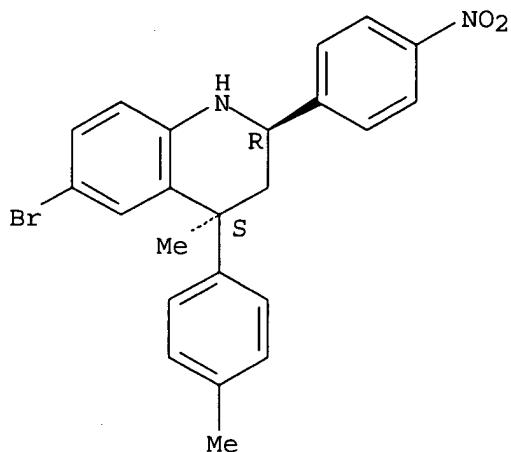
RN 635303-07-0 CAPLUS

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4/12/05

CN Quinoline, 6-bromo-1,2,3,4-tetrahydro-4-methyl-4-(4-methylphenyl)-2-(4-nitrophenyl)-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

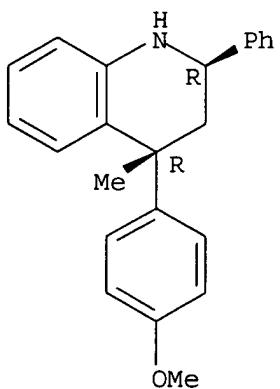
Relative stereochemistry.



RN 727653-41-0 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4-(4-methoxyphenyl)-4-methyl-2-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



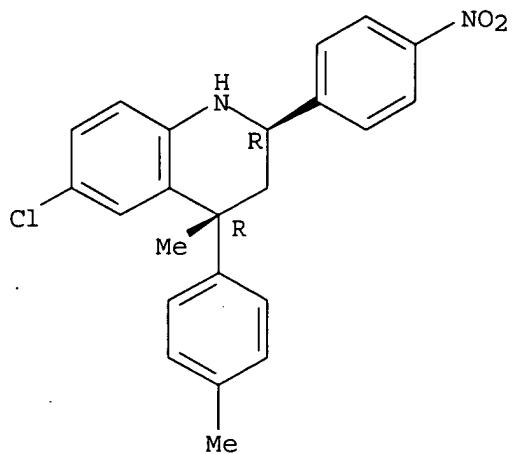
RN 727653-65-8 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-4-(4-methylphenyl)-2-(4-nitrophenyl)-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

10796396

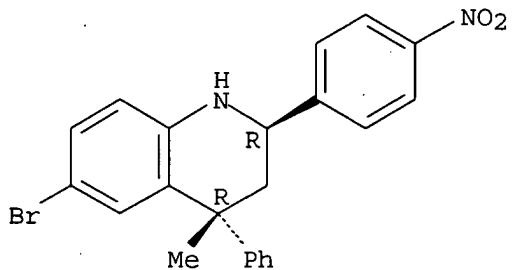
4/12/05



RN 727653-68-1 CAPLUS

CN Quinoline, 6-bromo-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

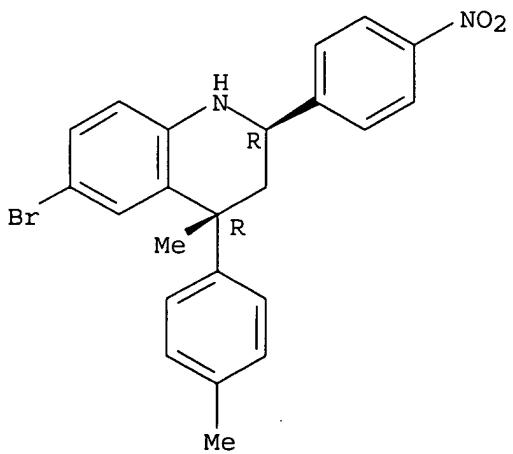
Relative stereochemistry.



RN 727653-73-8 CAPLUS

CN Quinoline, 6-bromo-1,2,3,4-tetrahydro-4-methyl-4-(4-methylphenyl)-2-(4-nitrophenyl)-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

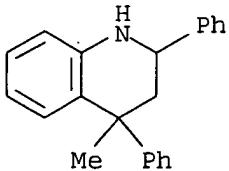


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RE.CNT 30 THERE ARE 30 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 4 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
AB A novel combined system of  $\text{Yb}(\text{OTf})_3$  with  $\text{TMSCl}$  or  $\text{TMSOTf}$  catalyzed an imino ene reaction. The reaction of  $\text{N-tosylbenzaldimine}$  with  $\alpha$ -methylstyrene proceeded smoothly to give  $\text{TsNHCHPhCH}_2\text{CPh:CH}_2$  in the presence of a catalytic amount of  $\text{Yb}(\text{OTf})_3$  and  $\text{TMSCl}$ . This catalytic system was successfully applied to the imino ene reactions of various aldimines with alkenes. This new imino ene reaction provides a unique method for the three-component coupling reaction of an aldehyde, tosylamide, and  $\alpha$ -methylstyrene in the presence of  $\text{Yb}(\text{OTf})_3$  and  $\text{TMSOTf}$ , to give the homoallylic amine.  
AN 2003:213238 CAPLUS  
DN 138:385135  
TI Imino Ene Reaction Catalyzed by Ytterbium(III) Triflate and  $\text{TMSCl}$  or  $\text{TMSOTf}$   
AU Yamanaka, Masamichi; Nishida, Atsushi; Nakagawa, Masako  
CS Graduate School of Pharmaceutical Sciences, Chiba University, Inage, Chiba, 263-8522, Japan  
SO Journal of Organic Chemistry (2003), 68(8), 3112-3120  
CODEN: JOCEAH; ISSN: 0022-3263  
PB American Chemical Society  
DT Journal  
LA English  
OS CASREACT 138:385135  
IT 30312-28-8P  
RL: SPN (Synthetic preparation); PREP (Preparation)  
(imino ene reaction catalyzed by ytterbium(III) triflate and  $\text{TMSCl}$  or  $\text{TMSOTf}$ )  
RN 30312-28-8 CAPLUS  
CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl- (9CI) (CA INDEX NAME)



RE.CNT 49 THERE ARE 49 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 5 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
AB Irradiation ( $\lambda > 345$  nm) of catalytic amts. of 2,4,6-triphenylpyrylium tetrafluoroborate (TPT) in a  $\text{CH}_2\text{Cl}_2$  solution of  $\text{N-aryl imines}$  and  $\alpha$ -methylstyrene or styrene produced the corresponding [4+2] cycloaddn. products, tetrahydroquinoline derivs., in good yield. The reaction was controlled by the relative oxidation potentials of the dienophile and the diene.  
AN 2002:880560 CAPLUS  
DN 138:287505  
TI Photosensitized Diels-Alder reactions of  $\text{N-arylimines}$ : synthesis of tetrahydroquinoline derivatives  
AU Zhang, Wei; Jia, Xiaodong; Yang, Li; Liu, Zhong-Li

10796396

4/12/05

CS National Laboratory of Applied Organic Chemistry, Lanzhou University,  
Gansu, 730000, Peop. Rep. China

SO Tetrahedron Letters (2002), 43(51), 9433-9436  
CODEN: TELEAY; ISSN: 0040-4039

PB Elsevier Science Ltd.

DT Journal

LA English

OS CASREACT 138:287505

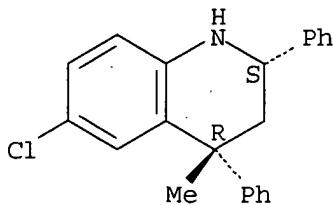
IT 184226-21-9P 258851-87-5P 258851-91-1P  
506407-49-4P 506407-50-7P 506407-51-8P  
506407-52-9P 506407-53-0P 506407-54-1P  
506407-60-9P 506407-61-0P 506407-62-1P  
506407-63-2P 506407-64-3P 506407-65-4P  
506407-66-5P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation of tetrahydroquinoline derivs. via photosensitized Diels-Alder  
reactions of N-arylimines and styrenes)

RN 184226-21-9 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-,  
(2R,4S)-rel- (9CI) (CA INDEX NAME)

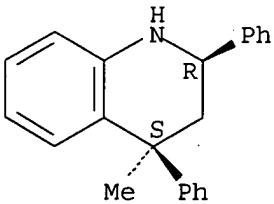
Relative stereochemistry.



RN 258851-87-5 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-, (2R,4S)-rel- (9CI)  
(CA INDEX NAME)

Relative stereochemistry.



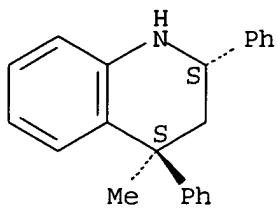
RN 258851-91-1 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-, (2R,4R)-rel- (9CI)  
(CA INDEX NAME)

Relative stereochemistry.

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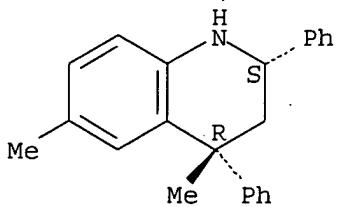
4/12/05



RN 506407-49-4 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2,4-diphenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

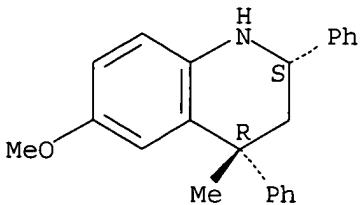
Relative stereochemistry.



RN 506407-50-7 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2,4-diphenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

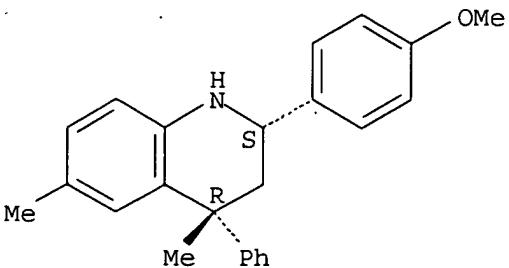
Relative stereochemistry.



RN 506407-51-8 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-2-(4-methoxyphenyl)-4,6-dimethyl-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 506407-52-9 CAPLUS

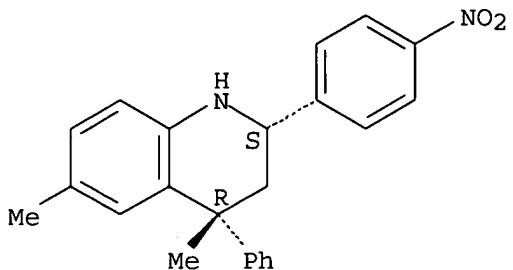
CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2-(4-nitrophenyl)-4-phenyl-,

10796396

4/12/05

(2R,4S)-rel- (9CI) (CA INDEX NAME)

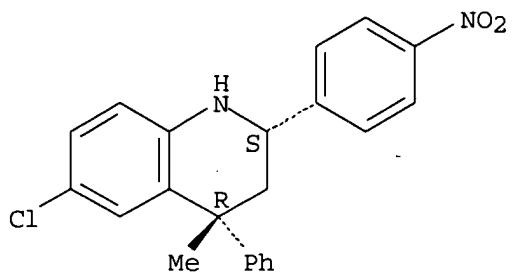
Relative stereochemistry.



RN 506407-53-0 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

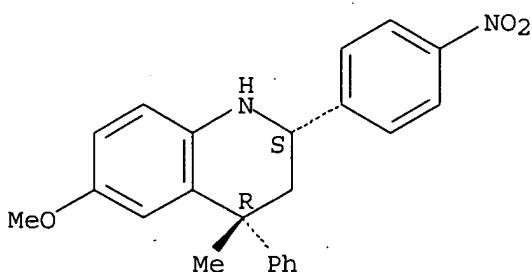
Relative stereochemistry.



RN 506407-54-1 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



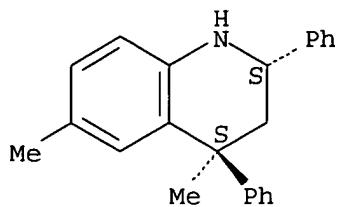
RN 506407-60-9 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2,4-diphenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

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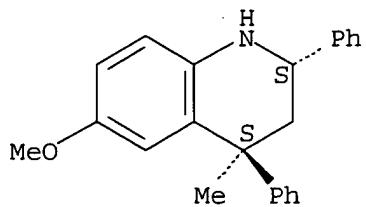
4/12/05



RN 506407-61-0 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2,4-diphenyl-,  
(2R,4R)-rel- (9CI) (CA INDEX NAME)

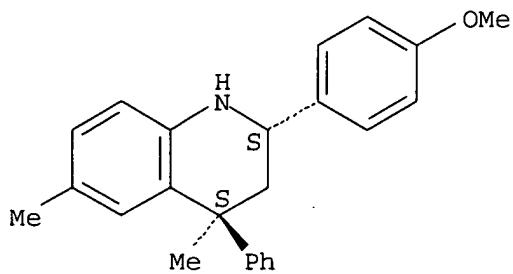
Relative stereochemistry.



RN 506407-62-1 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-2-(4-methoxyphenyl)-4,6-dimethyl-4-phenyl-,  
(2R,4R)-rel- (9CI) (CA INDEX NAME)

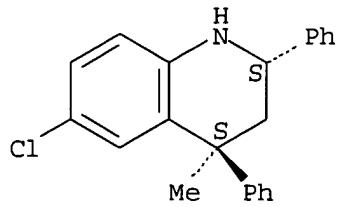
Relative stereochemistry.



RN 506407-63-2 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-,  
(2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 506407-64-3 CAPLUS

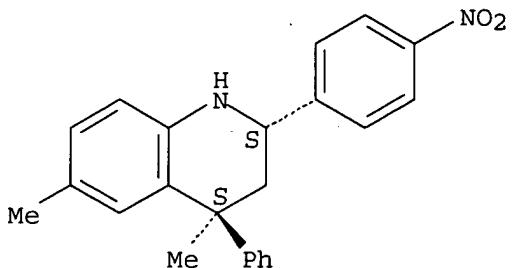
CN Quinoline, 1,2,3,4-tetrahydro-4,6-dimethyl-2-(4-nitrophenyl)-4-phenyl-,

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(2R,4R)-rel- (9CI) (CA INDEX NAME)

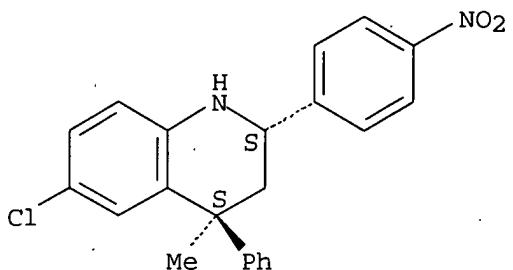
Relative stereochemistry.



RN 506407-65-4 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

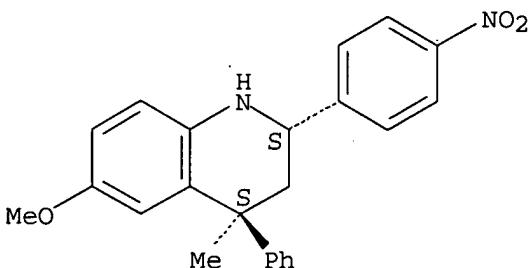
Relative stereochemistry.



RN 506407-66-5 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-6-methoxy-4-methyl-2-(4-nitrophenyl)-4-phenyl-, (2R,4R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RE.CNT 39 THERE ARE 39 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 6 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN

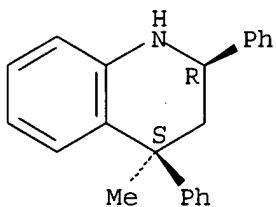
AB Ytterbium trifluoromethanesulfonate [Yb(OTf)<sub>3</sub>] catalyzed the imino ene reaction of N-tosyl aldimine with  $\alpha$ -methylstyrene to give a homoallylamine in moderate yield. Furthermore, addition of a catalytic amount of chlorotrimethylsilane (TMSCl) dramatically enhanced the imino ene reaction.

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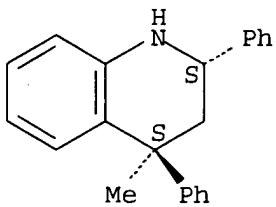
AN 1999:803639 CAPLUS  
DN 132:165964  
TI Ytterbium(III) Triflate/TMSCl: Efficient Catalyst for Imino Ene Reaction  
AU Yamanaka, Masamichi; Nishida, Atsushi; Nakagawa, Masako  
CS Faculty of Pharmaceutical Science, Chiba University, Inage-ku Chiba,  
263-8522, Japan  
SO Organic Letters (2000), 2(2), 159-161  
CODEN: ORLEF7; ISSN: 1523-7060  
PB American Chemical Society  
DT Journal  
LA English  
OS CASREACT 132:165964  
IT 258851-87-5P 258851-91-1P  
RL: SPN (Synthetic preparation); PREP (Preparation)  
(ytterbium(III) triflate/TMSCl catalyst system for the imino ene  
reaction of N-tosyl aldimine with  $\alpha$ -methylstyrene)  
RN 258851-87-5 CAPLUS  
CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-, (2R,4S)-rel- (9CI)  
(CA INDEX NAME)

Relative stereochemistry.



RN 258851-91-1 CAPLUS  
CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-, (2R,4R)-rel- (9CI)  
(CA INDEX NAME)

Relative stereochemistry.

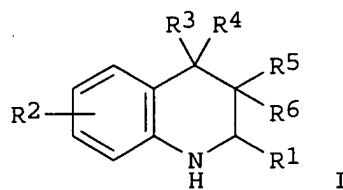


RE.CNT 35 THERE ARE 35 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 7 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
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4/12/05



AB A convenient method for library construction in liquid phase, which is based on lanthanide triflate ( $\text{Ln}(\text{OTf})_3$ )-catalyzed three-component reactions, has been developed. Equimolar amts. of each component, an aldehyde, an amine, and a silyl enol ether or an alkene, react smoothly in the presence of  $\text{Ln}(\text{OTf})_3$ , and the work-up and purification processes are performed by simply passing through a short column. The key is to use  $\text{Ln}(\text{OTf})_3$  as a Lewis acid catalyst, which is not decomposed during the work-up and purification steps,

and is easily separated from products by the simple procedure. According to this method, various high-quality  $\beta$ -amino ester and quinoline derivs.  $\text{R}_2\text{NHCHR}_1\text{CR}_3\text{R}_4\text{COR}_5$  ( $\text{R}_1 = \text{Ph}$ , 2-furyl,  $\text{PhCO}$ , etc.,  $\text{R}_2 = \text{Ph}$ ,  $\text{PhCH}_2$ , 4-MeOC<sub>6</sub>H<sub>4</sub>,  $\text{R}_3 = \text{Me}$ , H,  $\text{PhCH}_2\text{O}$ ,  $\text{R}_4 = \text{Me}$ , H,  $\text{OSiMe}_2\text{CMe}_3$ ,  $\text{R}_5 = \text{OMe}$ ,  $\text{SEt}$ ,  $\text{Ph}$ ) and I ( $\text{R}_1 = \text{Ph}$ , 4-ClC<sub>6</sub>H<sub>4</sub>,  $\alpha$ -naphthyl, etc.,  $\text{R}_2 = \text{H}$ , 4-Cl, cyclohexyl, etc.,  $\text{R}_3$ ,  $\text{R}_4$ ,  $\text{R}_5$ ,  $\text{R}_6$  = alkene substituent) have been prepared in parallel in large quantities.

AN 1998:243718 CAPLUS

DN 128:270525

TI A convenient method for library construction: parallel synthesis of  $\beta$ -amino ester and quinoline derivatives in liquid phase using  $\text{Ln}(\text{OTf})_3$ -catalyzed three-component reactions

AU Kobayashi, Shu; Komiyama, Susumu; Ishitani, Haruro

CS Department of Applied Chemistry, Faculty of Science, Science University of Tokyo, Tokyo, 162, Japan

SO Biotechnology and Bioengineering (1998), 61(1), 23-31

CODEN: BIBIAU; ISSN: 0006-3592

PB John Wiley & Sons, Inc.

DT Journal

LA English

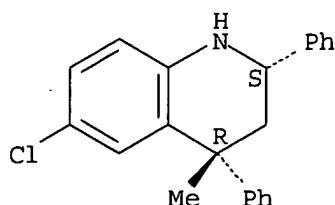
IT 184226-21-9P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(parallel synthesis of  $\beta$ -amino ester and quinoline derivs. using  $\text{Ln}(\text{OTf})_3$ -catalyzed three-component reactions)

RN 184226-21-9 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-, (2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



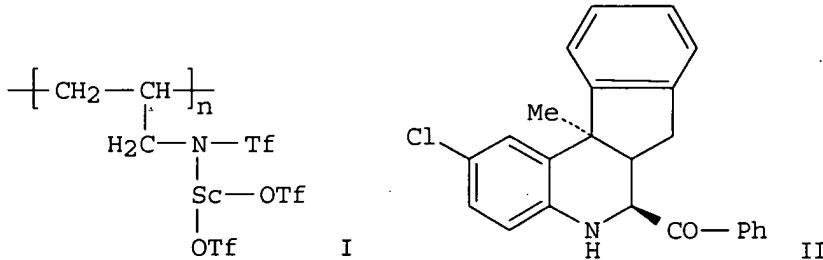
RE.CNT 59 THERE ARE 59 CITED REFERENCES AVAILABLE FOR THIS RECORD

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4/12/05

ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 8 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN  
GI



AB The poly(allylamine)-supported scandium catalyst I was prepared Quinolines were synthesized from anilines, arylaldehydes and alkenes in the presence of I. For example, the condensation of  $\alpha$ -oxobenzeneacetaldehyde with 4-chloroaniline and 3-methyl-1H-indene gave II in 99% yield as a single diastereomer.

AN 1996:619167 CAPLUS

DN 126:31256

TI A New Methodology for Combinatorial Synthesis. Preparation of Diverse Quinoline Derivatives Using a Novel Polymer-Supported Scandium Catalyst

AU Kobayashi, Shu; Nagayama, Satoshi

CS Faculty of Science, Science University of Tokyo (SUT), Tokyo, 162, Japan

SO Journal of the American Chemical Society (1996), 118(37), 8977-8978

CODEN: JACSAT; ISSN: 0002-7863

PB American Chemical Society

DT Journal

LA English

OS CASREACT 126:31256

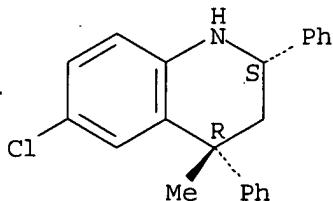
IT 184226-21-9P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation of quinoline combinatorial library via poly(allylamine)-  
supported scandium catalyzed condensation reaction)

RN 184226-21-9 CAPLUS

CN Quinoline, 6-chloro-1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl-,  
(2R,4S)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



L4 ANSWER 9 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN

GI For diagram(s), see printed CA Issue.

AB Tetrahydroquinolines I (R = Me, CHMe<sub>2</sub>, Pr, n-C<sub>5</sub>H<sub>11</sub>; R<sub>1</sub> = H, 6-Me, 6-OMe,

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8-Me, 5,8-Me<sub>2</sub>) were prepared in 29-76% yield by condensing R<sub>1</sub>C<sub>6</sub>N<sub>4</sub>NH<sub>2</sub> with PhCMe:CH<sub>2</sub> and RCHO. 2,4-Dimethyl-4-vinyl-1,2,3,4-tetrahydroquinoline was similarly prepared from isoprene and 4-dimethyl derivs. of I from styrene.

AN 1973:515454 CAPLUS

DN 79:115454

TI 1,2,3,4-Tetrahydroquinolines

PA Chemische Werke Huels A.-G.

SO Fr., 15 pp.

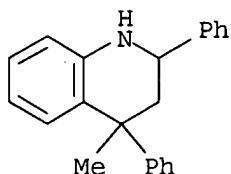
CODEN: FRXXAK

DT Patent

LA French

FAN CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	FR 1605038	A	19720828	FR 1968-1605038	19681211
	DE 1770170	A	19701223	DE 1967-1770170	19680410
PRAI	DE 1967-1770170	A	19680410		
IT	30312-28-8P			RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of)	
RN	30312-28-8	CAPLUS			
CN	Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl- (9CI)			(CA INDEX NAME)	



L4 ANSWER 10 OF 10 CAPLUS COPYRIGHT 2005 ACS on STN

GI For diagram(s), see printed CA Issue.

AB RC<sub>6</sub>H<sub>4</sub>NH<sub>2</sub> (I) (where R = H, p-Me, p-MeO, p-Cl, o-NO<sub>2</sub>, or p-NO<sub>2</sub>), R<sub>1</sub>CHO (II) (where R<sub>1</sub> = Me, iso-Pr, C<sub>5</sub>H<sub>11</sub>, or Ph), and R<sub>2</sub>C(R<sub>3</sub>):CH<sub>2</sub> (III) (where R<sub>2</sub> = H or Me, R<sub>3</sub> = Ph or CH:CH<sub>2</sub>) in HOAc and concentrated H<sub>2</sub>SO<sub>4</sub> gave by cyclization

the corresponding title compds. (IV). II (R<sub>1</sub> = H) gave in this reaction predominantly 1-R<sub>2</sub>,1-R<sub>3</sub>,7-R<sub>2</sub>,7-R<sub>3</sub>-tetrasubstituted-2,3,6,7-tetrahydro-1H,5H-benzo[ij]quinolizines (V).

AN 1971:42266 CAPLUS

DN 74:42266

TI Synthesis of 1,2,3,4-tetrahydroquinolines

AU Hesse, Karl D.

CS Forschungslab., Chem. Werke Huels A.-G., Marl, Fed. Rep. Ger.

SO Justus Liebigs Annalen der Chemie (1970), 741, 117-23

CODEN: JLACBF; ISSN: 0075-4617

DT Journal

LA German

OS CASREACT 74:42266

IT 30312-28-8P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation of)

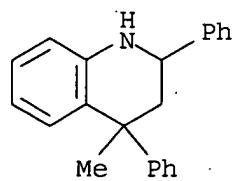
RN 30312-28-8 CAPLUS

CN Quinoline, 1,2,3,4-tetrahydro-4-methyl-2,4-diphenyl- (9CI) (CA INDEX

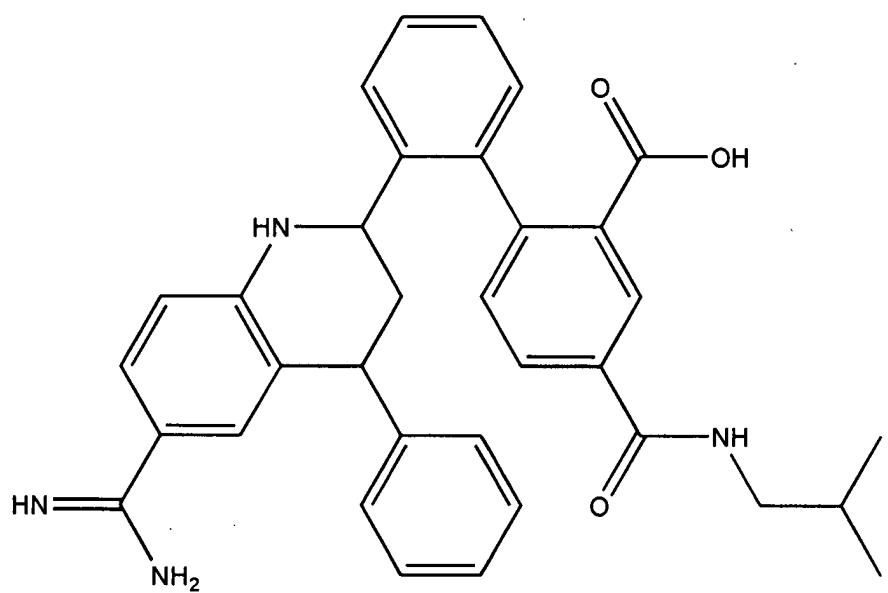
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NAME)



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2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-isobutylcarbamoyl-biphenyl-2-carboxylic acid

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L8 FILE 'CAPLUS' ENTERED AT 16:42:36 ON 12 APR 2005  
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S L8

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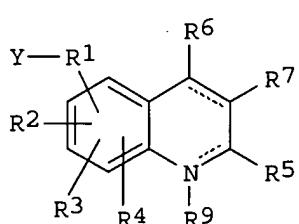
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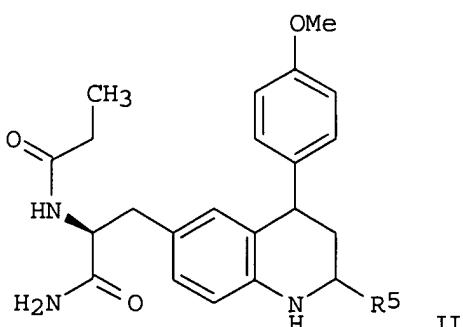
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L16 1 L15 NOT L12

=> d abs fbib hitstr

L16 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN  
GI



I



II

AB The invention relates to novel 4-substituted quinoline derivs. I, their salts, and combinatorial libraries containing mixts. of two or more such compds. [wherein R1 = bond, (un)substituted alk(en/yn)ylene, cycloalk(en)ylene, phenylene, naphthylene, heterocycle, heteroaryl, amino,

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CH<sub>2</sub>CONH, (CH<sub>2</sub>)<sub>p</sub>Ar(CH<sub>2</sub>)<sub>q</sub>, etc.; p, q = 0-6 but both cannot be 0; Ar = (un)substituted Ph or heteroaryl; R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub> = H, halo, (un)protected OH, cyano, NO<sub>2</sub>, (un)substituted alk(en/yn)yl, alkoxy, cycloalk(en)yl, heterocyclyl, phenylalkyl, Ph, naphthyl, etc.; R<sub>5</sub> = H, (un)substituted alk(en/yn)yl, cycloalk(en)yl, Ph, naphthyl, phenylalkyl, (un)protected CO<sub>2</sub>H, acyl, heterocyclyl, etc.; R<sub>6</sub> = H, (un)substituted Ph, naphthyl, 2-oxopyrrolidin-1-yl and higher homologs, (un)substituted NHCHO; R<sub>7</sub> = H, (un)substituted alkyl; Y = CO<sub>2</sub>H, OH, SH, NHR<sub>8</sub>, CONHR<sub>8</sub>, CH<sub>2</sub>OH, CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>NHR<sub>8</sub>; R<sub>8</sub> = H, (un)substituted alkyl, or functionalized resin; R<sub>9</sub> = H, (un)substituted alkyl, phenylalkyl, acyl, PhSO<sub>2</sub>, alkylsulfonyl, alkylaminocarbonyl, or PhNHCO, or is absent; dotted lines = optional pi bonds]. The invention also relates to the generation of such libraries. In 12 examples, libraries of I ranging in size from 2380 to 39,440 compds. were prepared as mixed sublibraries. Data for control compds. (samples of individually known intermediates and products, cleaved from simultaneously processed control resins) are given for some examples. Both quinoline and tetrahydroquinoline libraries were prepared. For instance, tea-bags of MBHA resin were each coupled with L- or D-N-BOC-p-nitrophenylalanine, the BOC groups were removed from both, and the amino groups were each acylated with 170 carboxylic acids. The acylated, resin-bound products were mixed and reduced at the nitro group, and the amine product mixts. were condensed with 58 different aldehydes and cyclized with 4-methoxystyrene. Cleavage of the resin-bound products with HF gave mixed sublibraries of I. Individual control samples of products, such as II [R<sub>5</sub> = 1-naphthyl, 2,3-difluorophenyl, cyclohexyl, etc.], were obtained by reactions of pure, resin-bound L-N-propanoyl-p-aminophenylalanine control samples with individual aldehydes and 4-methoxystyrene. Potential applications of I (no data) may include use as antibacterials, NMDA antagonists, or analgesics.

AN 1998:543220 CAPLUS  
DN 129:175563  
TI 4-Substituted quinoline derivatives and 4-substituted quinoline combinatorial libraries

IN Hayes, Thomas K.; Forood, Behrouz; Kiely, John S.  
PA Trega Biosciences, Inc., USA  
SO PCT Int. Appl., 124 pp.

CODEN: PIXXD2

DT Patent  
LA English

FAN. CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9834115	A1	19980806	WO 1997-US22391	19971205
	W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
CA	2279977	AA	19980806	US 1997-795392 CA 1997-2279977 US 1997-795392 WO 1997-US22391	A 19970204 19971205 A 19970204 W 19971205
AU	9881919	A1	19980825	AU 1998-81919 US 1997-795392 WO 1997-US22391	19971205 A 19970204 W 19971205

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EP 977989	A1 20000209	EP 1997-949775	19971205
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI			
		US 1997-795392	A 19970204
		WO 1997-US22391	W 19971205
US 6262269	B1 20010717	US 1998-17785	19980203
		US 1997-126414P	P 19970204
		US 1997-795392	A 19970204
US 6388081	B1 20020514	US 1999-376670	19990816
		US 1997-126414P	P 19970204
		US 1998-17785	A3 19980203

OS MARPAT 129:175563

IT 211492-55-6P 211492-67-0P 211493-20-8P

211493-65-1P 211493-73-1P 211494-06-3P

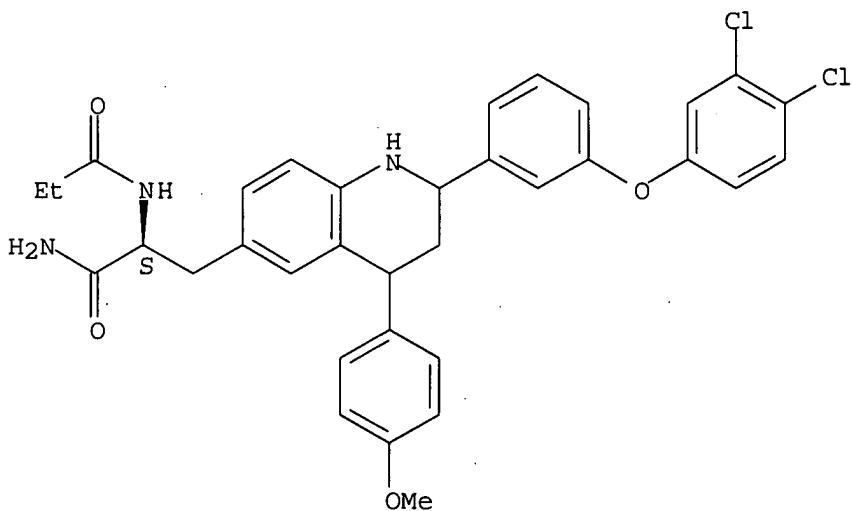
211494-15-4P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(resin-cleavage control product; preparation of 4-substituted quinoline  
derivs. and combinatorial libraries)

RN 211492-55-6 CAPLUS

CN 6-Quinolinepropanamide, 2-[3-(3,4-dichlorophenoxy)phenyl]-1,2,3,4-tetrahydro-4-(4-methoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]-, ( $\alpha$ S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



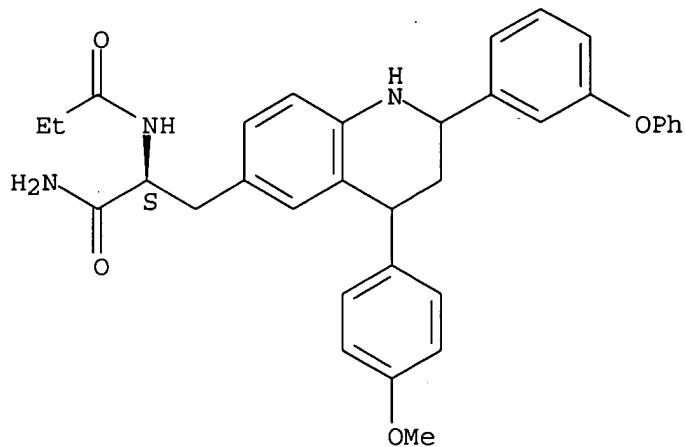
RN 211492-67-0 CAPLUS

CN 6-Quinolinepropanamide, 1,2,3,4-tetrahydro-4-(4-methoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]-2-(3-phenoxyphenyl)-, ( $\alpha$ S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

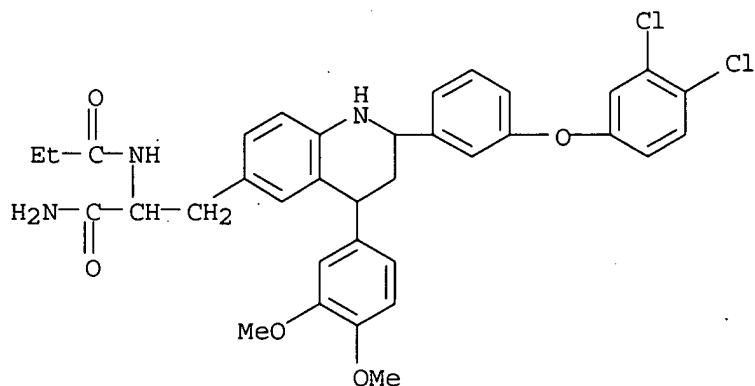
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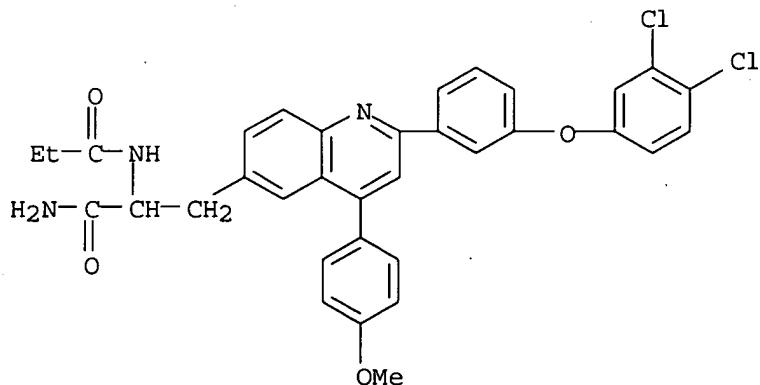
RN 211493-20-8 CAPLUS

CN 6-Quinolinepropanamide, 2-[3-(3,4-dichlorophenoxy)phenyl]-4-(3,4-dimethoxyphenyl)-1,2,3,4-tetrahydro- $\alpha$ -[(1-oxopropyl)amino]- (9CI)  
(CA INDEX NAME)



RN 211493-65-1 CAPLUS

CN 6-Quinolinepropanamide, 2-[3-(3,4-dichlorophenoxy)phenyl]-4-(4-methoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]- (9CI) (CA INDEX NAME)

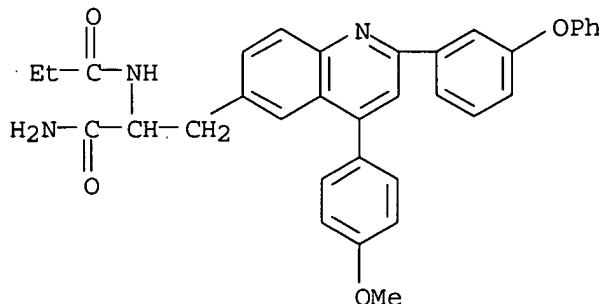


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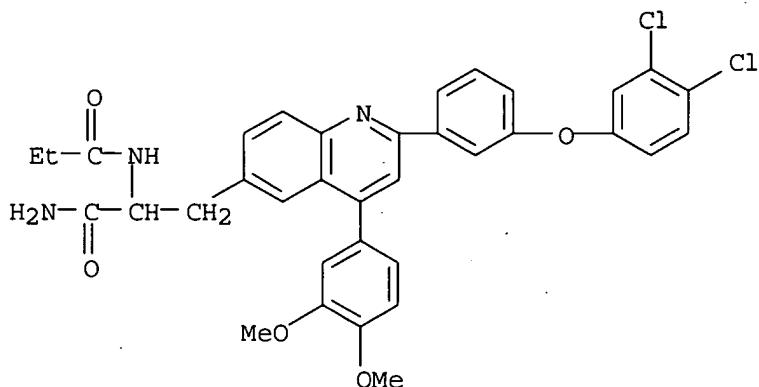
RN 211493-73-1 CAPLUS

CN 6-Quinolinepropanamide, 4-(4-methoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]-2-(3-phenoxyphenyl)- (9CI) (CA INDEX NAME)



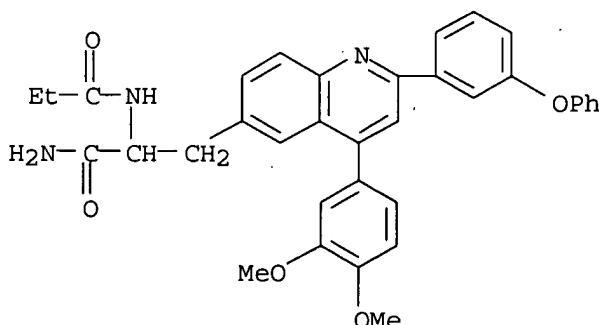
RN 211494-06-3 CAPLUS

CN 6-Quinolinepropanamide, 2-[3-(3,4-dichlorophenoxy)phenyl]-4-(3,4-dimethoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]- (9CI) (CA INDEX NAME)



RN 211494-15-4 CAPLUS

CN 6-Quinolinepropanamide, 4-(3,4-dimethoxyphenyl)- $\alpha$ -[(1-oxopropyl)amino]-2-(3-phenoxyphenyl)- (9CI) (CA INDEX NAME)



RE.CNT 4

THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

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=> file registry		
COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
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\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
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\*\*\*\*\*

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more  
information enter HELP PROP at an arrow prompt in the file or refer  
to the file summary sheet on the web at:  
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L17 STRUCTURE UPLOADED

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L17 HAS NO ANSWERS  
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\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SCREEN SEARCH COMPLETED - 70 TO ITERATE

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FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
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STRUCTURE FILE UPDATES: 11 APR 2005 HIGHEST RN 848290-51-7  
DICTIONARY FILE UPDATES: 11 APR 2005 HIGHEST RN 848290-51-7

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when  
conducting SmartSELECT searches.

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\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
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Crossover limits have been increased. See HELP CROSSOVER for details.

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<http://www.cas.org/ONLINE/DBSS/registryss.html>

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L20 STRUCTURE UPLOADED

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100.0% PROCESSED 21771 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

L21 0 SEA SSS FUL L20

10796396

The Examiner is requested to consider the foregoing information in relation to this application and indicate that each reference was considered by returning a copy of the initialed PTO 1449 form(s).

Respectfully submitted,

Bristol-Myers Squibb Company  
Patent Department  
P.O. Box 4000  
Princeton, NJ 08543-4000  
(609) 252-3791

  
\_\_\_\_\_  
Jing S. Belfield  
Agent for Applicants  
Reg. No. 45,914

Date: May 25, 2004